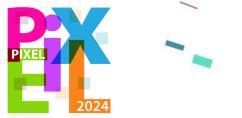
ELEVENTH INTERNATIONAL WORKSHOP ON SEMICONDUCTOR PIXEL DETECTORS FOR PARTICLES AND IMAGING, PIXEL2024



UNIVERSITY OF STRASBOURG, 18-22 NOVEMBER 2024

REALISTIC MONTE CARLO SIMULATION OF SILICON PARTICLE DETECTORS FOR TIMING & TRACKING WITH GARFIELD++

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## OUTLINE

- Introduction about Garfield++
- Simulation of solid-state detectors
  - electric field setup
  - signal formation
  - charge multiplication
  - read-out electronics
- Case-study I: monolithic CMOS detectors for timing
- Case-study II: monolithic CMOS detectors for 4D particle tracking
- Final Comments & Acknowledgments

"Garfield++ is an object-oriented toolkit for the detailed simulation of particle detectors based on ionization measurement in gases or semiconductors"

- **Garfield**++ is an evolution of the previous <u>Garfield</u> tool (*Rob Veenhof*, 1984) with which shares functionality, and from which differentiates by some point:
  - treatment of charge carrier transport
  - user interface (ROOT-based)
- Its main elements are:
  - ionization pattern: *relativistic particles*, *X-ray* absorption, (low-energy) *ions* or others
  - **transport** properties: transport and avalanche in *gas mixtures* or *semiconductors*
  - electric field calculation: different options available ...

#### ➤ Primary Ionization

- <u>relativistic particles</u>: <u>HEED</u> (High-Energy ElectroDynamics) program, extension of the <u>PAI</u> (Photo-Absorption Ionization) model
- <u>X-ray photoabsorption</u>: interface with HEED
- <u>Ion tracks</u>: imported results from <u>SRIM</u><sup>†</sup> or <u>TRIM</u><sup>‡</sup>. The first simulates the energy loss of ions in matter, the latter is a Monte Carlo extension that, in addition, accounts for individual ion trajectories
- Other: ionization/avalanche in gases (<u>Magboltz</u>), interface with GEANT4, ...

<sup>&</sup>lt;sup>†</sup>Stopping Range of Ions in Matter; <sup>‡</sup>TRansport of Ions in Matter

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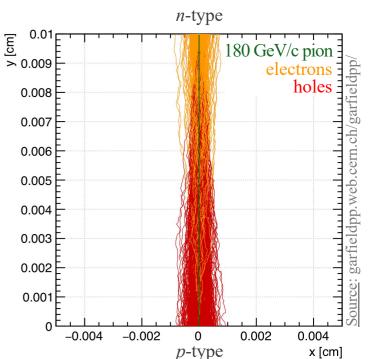
#### **INTRODUCTION ABOUT GARFIELD++**

#### ➤ Charge Transport

To find the **drift lines** of **individual electrons** and **holes**, the first-order equation of motion has to be solved.

Garfield++ offers (at least) two different strategies:

- <u>Runge-Kutta-Fehlberg</u> (RKF) integration, proceeding by iteratively estimating the charge position, drift velocity and time step
- <u>Monte Carlo</u> (MC) integration, based on macroscopic transport parameters
- ...plus even more accurate methods: see few slides ahead



- ➤ Charge Transport: the Monte Carlo integration
- a step of length  $\Delta s = v_d \Delta t$  in the direction of the drift velocity  $v_d$ , relative to the local electric and magnetic field, is calculated. <u>Time-step</u>  $\Delta t$  or <u>distance-step</u>  $\Delta s$  must be specified by the user
- a random diffusion step is sampled from three uncorrelated Gaussian distributions: one with standard deviation  $\sigma_{//} = D_{//} \sqrt{\Delta s}$  for the component parallel to the drift velocity and two with standard deviations  $\sigma_{T} = D_{T} \sqrt{\Delta s}$  for the transverse components
- the **location** is updated by vectorially adding the two steps

#### ➤ Electric field

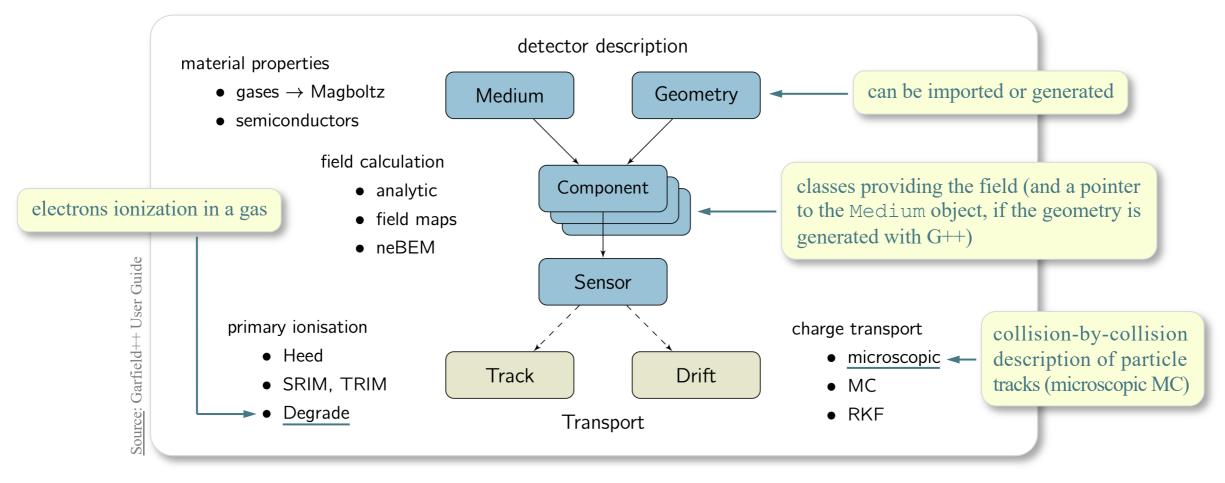
The field maps can be **imported** from <u>text files</u> or <u>external FEM tools</u> like

- Ansys
- Synopsys<sup>®</sup> Sentaurus TCAD
- Elmer
- <u>CST Studio</u>
- COMSOL Multiphysics

as well as **defined analytically** (only simple structures) or **directly computed**\*

<sup>\*</sup> through an interface with the **neBEM** (nearly exact Boundary Element Method) solver, or applying a **thin–wire limit** solution

#### **Overview of the Simulation Flow**



Typical simulation-based design process of **silicon detectors** for **HEP applications**:

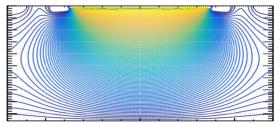
- 1. geometry definition with TCAD tools (2D or 3D)<sup>1</sup>
- 2. definition of **active electrodes**<sup>2</sup> for the signal acquisition
- 3. **exporting field** and **potential** maps (<u>no need to redefine the</u> <u>mesh in Garfield++</u>)
- 4. calculation of weighting field and weighting potential
- 5. **MC simulation** of **relativistic particles** crossing the detector (HEED model)
- 6. possible addition of readout electronic noise

TCAD

+ 5 geometry, mesh, doping  $\Rightarrow$  electric field, potential



weighting field, weighting potential  $\Rightarrow$  signals



<sup>&</sup>lt;sup>1</sup> to implement arrays, the user can define a periodicity of the geometry

<sup>&</sup>lt;sup>2</sup> to generate a weighting field/potential, one has to induce a small perturbation  $\delta V$  (~1% of  $V_{\text{bias}}$ ) on the collection electrode

To generate signals, once the weighting field has been generated, one has to set:

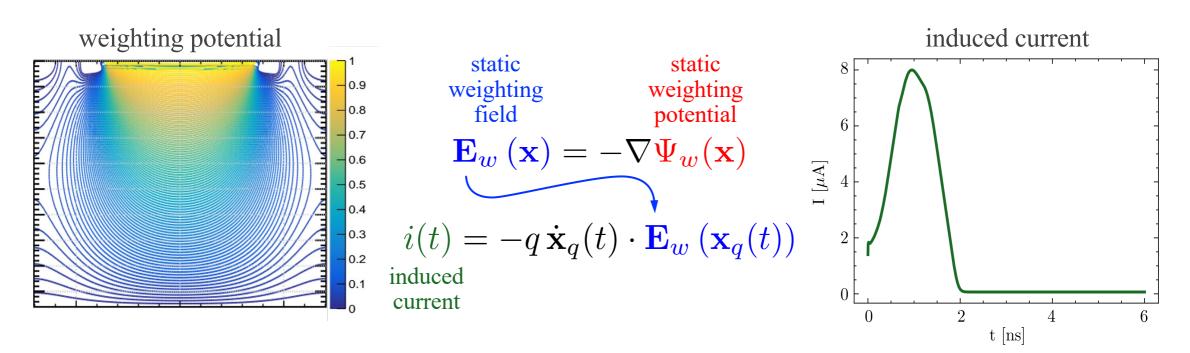
- the incident charged particle type:  $e^-$ ,  $e^+$ ,  $\mu^-$ ,  $\mu^+$ ,  $\pi^-$ ,  $\pi^+$ ,  $K^-$ ,  $K^+$ , p,  $\overline{p}$ , d
- its **momentum**
- the **trajectory**:
  - initial position and time vector (x0, y0, z0, t0)
  - direction vector (dx0, dy0, dz0)

Besides selecting the **distance-step**  $\Delta s$  (or **time-step**  $\Delta t$ ) for the charge transport process, we also need to set the **granularity** of the signal acquisition, through:

- initial time
- <u>final time</u>
- number of bins

Now we are ready to generate a signal

- apply a potential step  $\delta V$  to the readout electrode:  $\Psi_w = \frac{\Psi(V + \delta V) \Psi(V)}{\delta V}$
- compute the induced current through the Ramo-Shockley theorem



In presence of **internal gain**, there are four options for the **ionization coefficients** 

- **Grant:**  $\alpha_{\text{Gra}}(E) = A_{\text{Gra}} \exp\left(-\frac{B_{\text{Gra}}}{E}\right)$
- Massey:  $\alpha_{\text{Mas}}(E) = A_{\text{Mas}} \exp\left(-\frac{B_{\text{Mas}}(T)}{E}\right)$ , with  $B_{\text{Mas}}(T) = C_{\text{Mas}} + D_{\text{Mas}} \cdot T$
- van Overstraeten de Man:  $\alpha_{vOv}(E) = \gamma A_{vOv} \exp\left(-\gamma \frac{B_{vOv}}{E}\right)$

• **Okuto – Crowell:**  $\alpha_{\text{Oku}}(E) = A_{\text{Oku}} \left( 1 + (T - 300) C_{\text{Oku}} \right) E \exp \left( - \left( \frac{B_{\text{Oku}} \left( 1 + (T - 300) D_{\text{Oku}} \right)}{E} \right)^2 \right)$ 

The default model is van Overstraeten – de Man, with coefficients

 $\begin{aligned} A_{\rm vOv,n} &= 7.030 \cdot 10^5 \,{\rm cm}^{-1} & B_{\rm vOv,n} = 1.231 \cdot 10^6 \,{\rm V/cm} \\ A_{\rm vOv,p} &= 1.582 \cdot 10^6 \,{\rm cm}^{-1} \,\,(E < 400 \,{\rm kV/cm}) & B_{\rm vOv,p} = 2.036 \cdot 10^6 \,{\rm V/cm} \,\,(E < 400 \,{\rm kV/cm}) & \gamma = \frac{\tanh\left(\frac{\hbar\omega_{\rm op}}{2k_{\rm B}300{\rm K}}\right)}{\tanh\left(\frac{\hbar\omega_{\rm op}}{2k_{\rm B}T}\right)} \\ A_{\rm vOv,p} &= 6.710 \cdot 10^5 \,{\rm cm}^{-1} \,\,(E > 400 \,{\rm kV/cm}) & B_{\rm vOv,p} = 1.693 \cdot 10^6 \,{\rm V/cm} \,\,(E > 400 \,{\rm kV/cm}) & \gamma = \frac{\tanh\left(\frac{\hbar\omega_{\rm op}}{2k_{\rm B}T}\right)}{\tanh\left(\frac{\hbar\omega_{\rm op}}{2k_{\rm B}T}\right)} \end{aligned}$ 

Garfield++ allows the modeling of the **front-end electronics** by convoluting the induced current signal with a **transfer function** (delta response function)

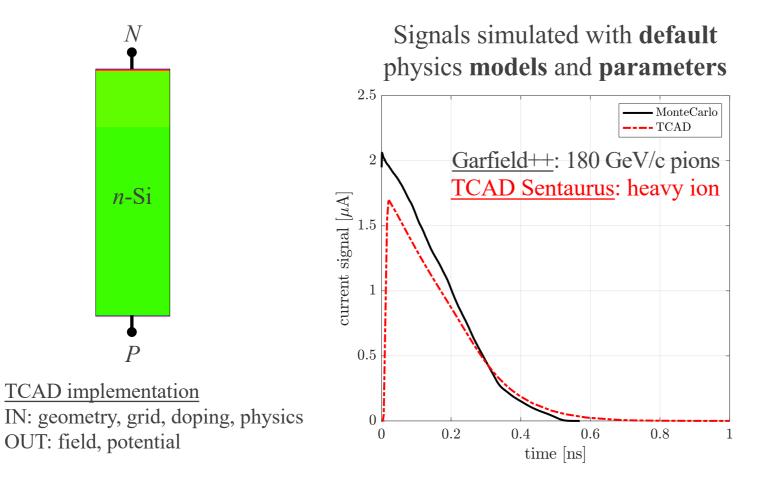
The transfer function can be:

- a user-specific <u>function</u>
- a <u>table</u>
- a pre-implemented <u>analytic model</u>, like the following *n*-stage shaper

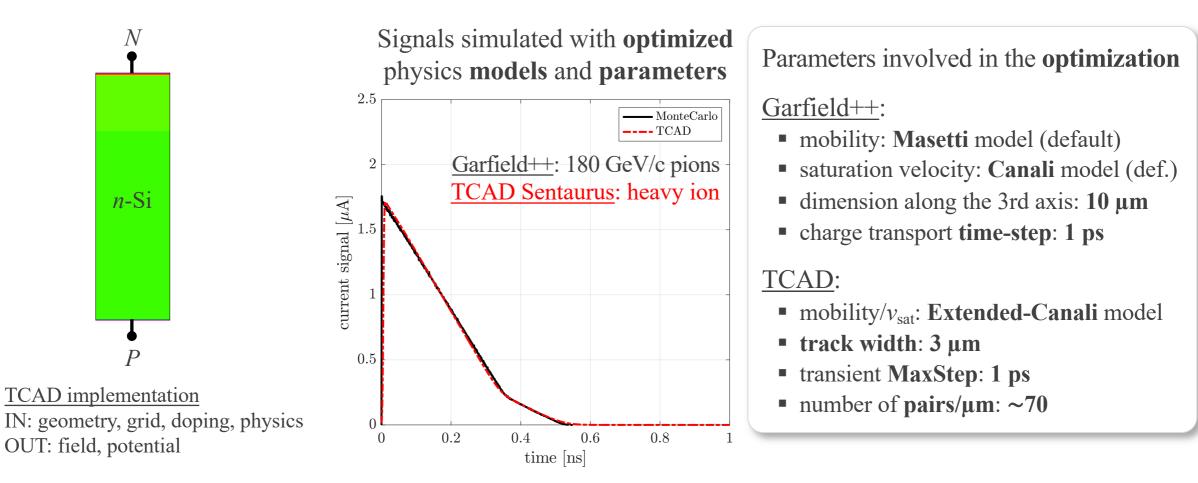
$$f(t) = g \exp(n) \left(\frac{t}{t_p}\right)^n \exp(-t/\tau), \text{ with } t_p = n\tau$$

Furthermore, also the **noise** can be added, reproducing a given equivalent-noise charge at the amplifier output

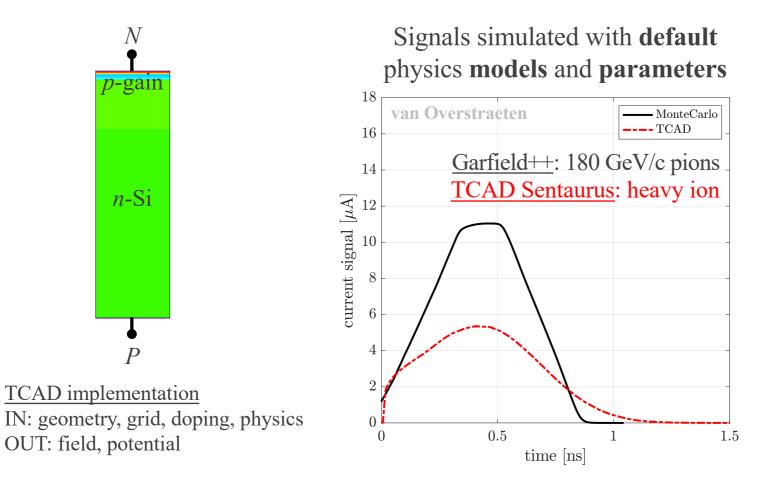
#### Example n.1: the *p-i-n* diode



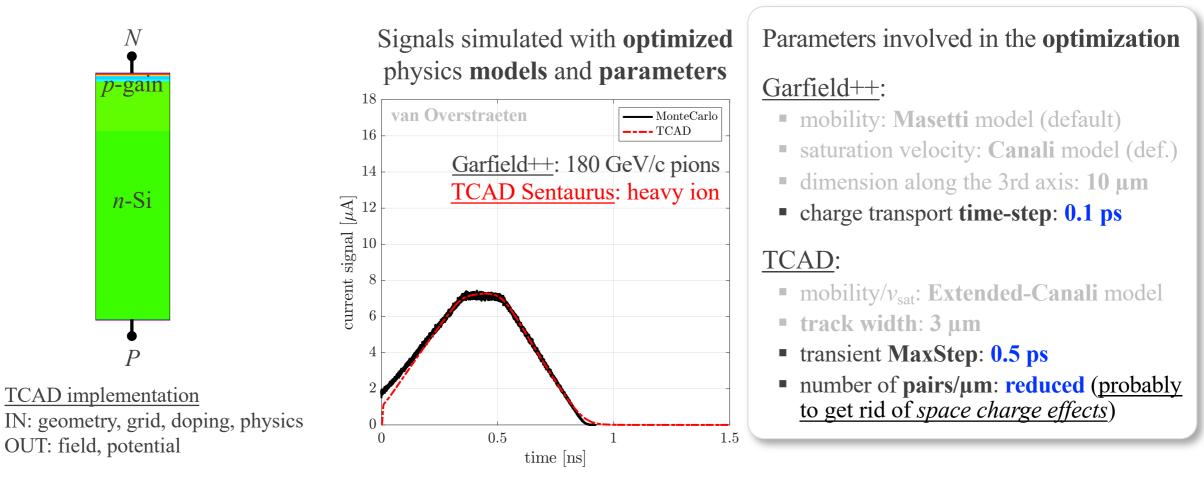
#### Example n.1: the *p-i-n* diode



#### Example n.2: the LGAD

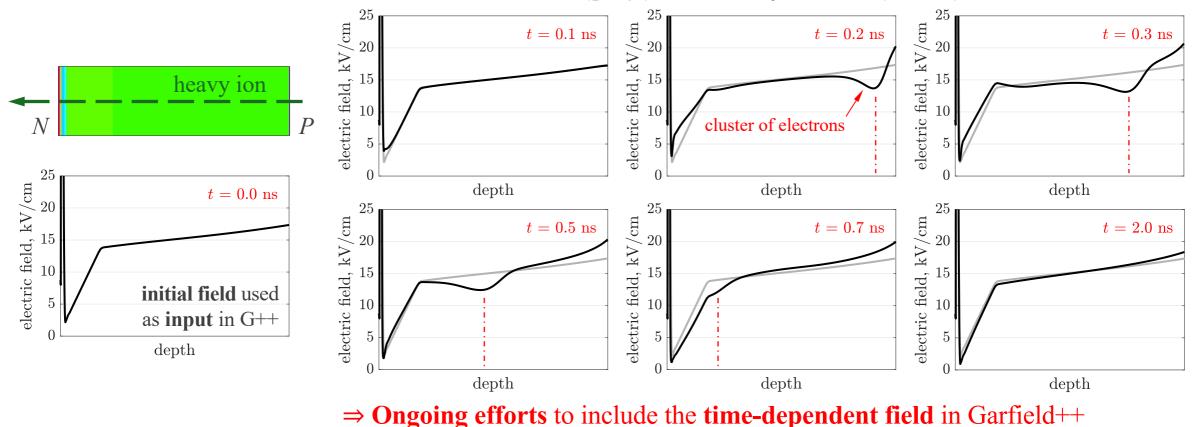


#### Example n.2: the LGAD

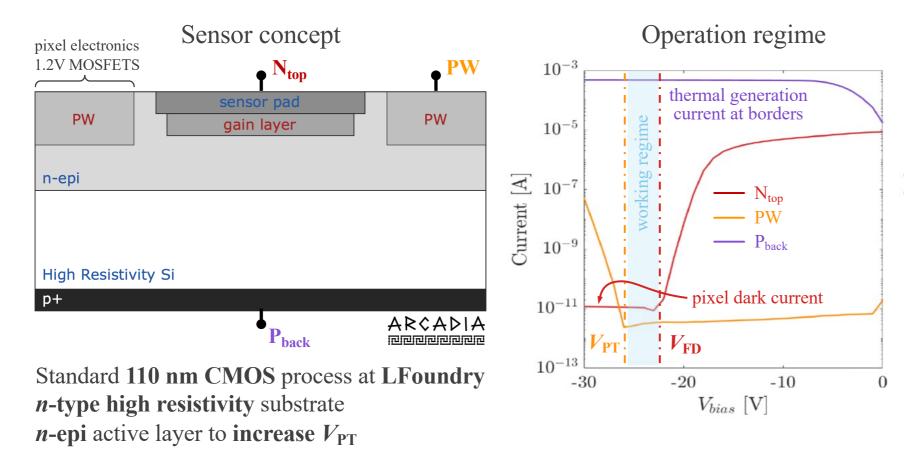


#### Example n.2: the LGAD

**Initial** *t* = 0 (*grey*) versus **dynamic** (*black*) **electric field** 



#### ARCADIA project (INFN) towards the development of MAPS with gain

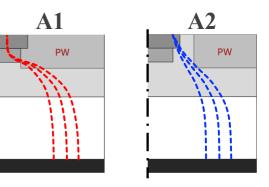


Main constraints

- **full-depletion** condition
- punch-through driven by the backside bias
- edge breakdown (due to gain) induced by the topside voltage

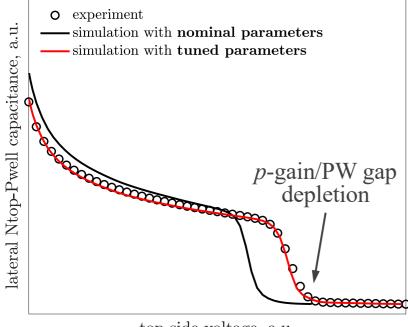
Two border layouts

- A1 uniform multiplication
- A2 (std.) uniform response



To have **realistic predictions** of the sensor figures-of-merit, the **technology parameters** have been previously tuned on **electrical measurements** 

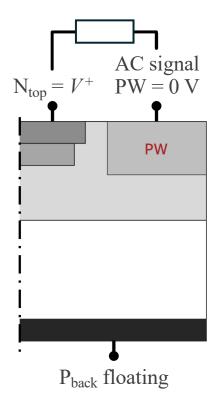
Lateral pad capacitance



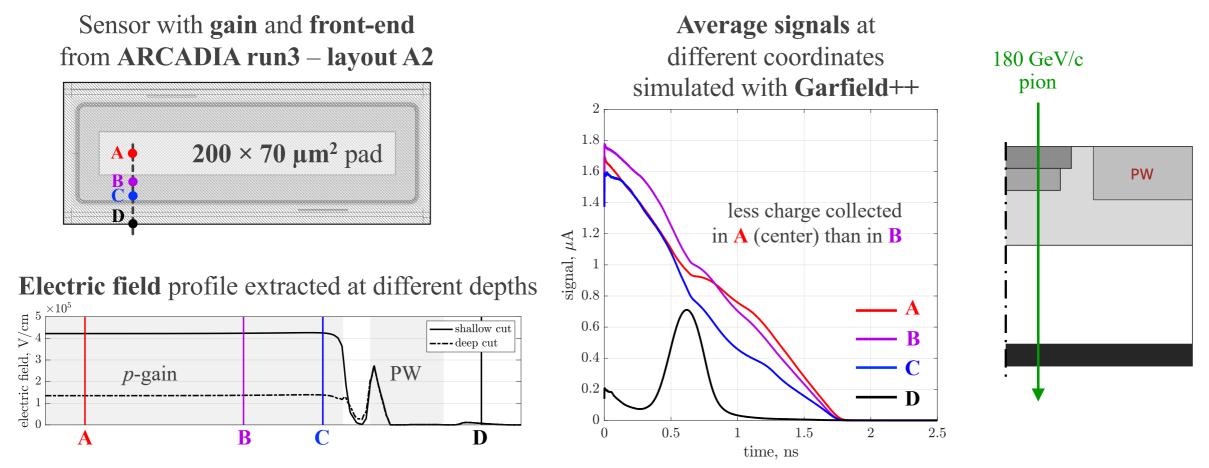
top-side voltage, a.u.

Parameters involved in the tuning:

- implants lateral diffusion (*p*-gain, n<sup>+</sup>, p-wells)
- *p*-gain **implantation energy** and **dose**
- *n*-epi doping concentration



#### Particle scan to investigate the impact of borders on charge collection

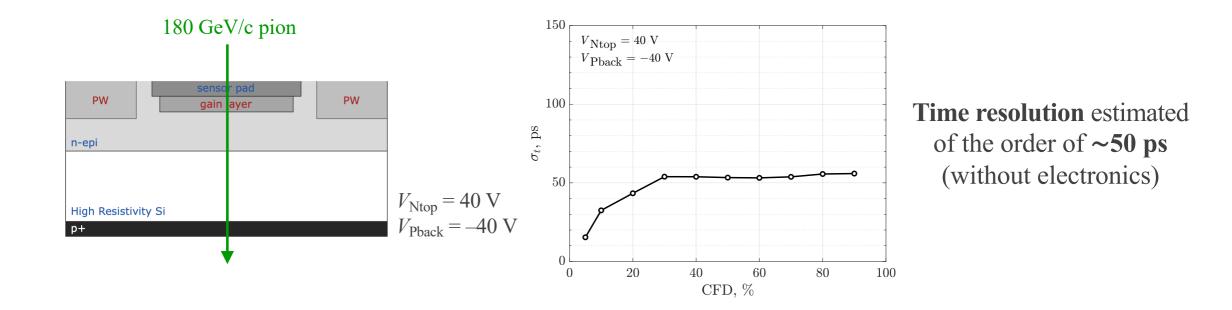


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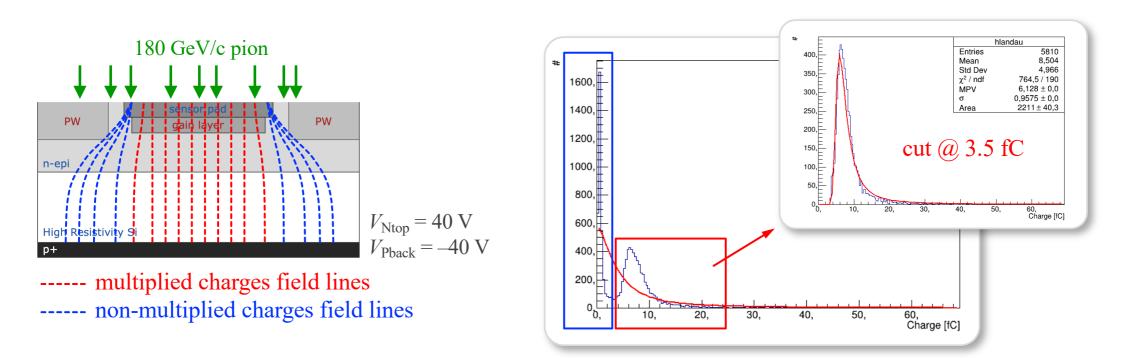
Further numerical investigations on the response uniformity:

• Particle injection in the center of the active area



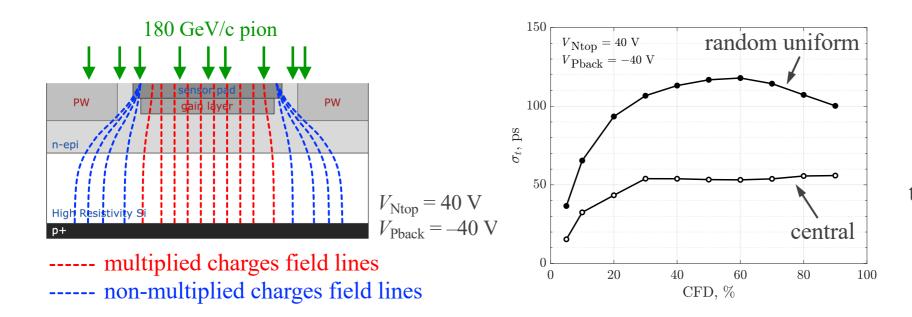
Further numerical investigations on the response uniformity:

Particle injection according to a random uniform distribution



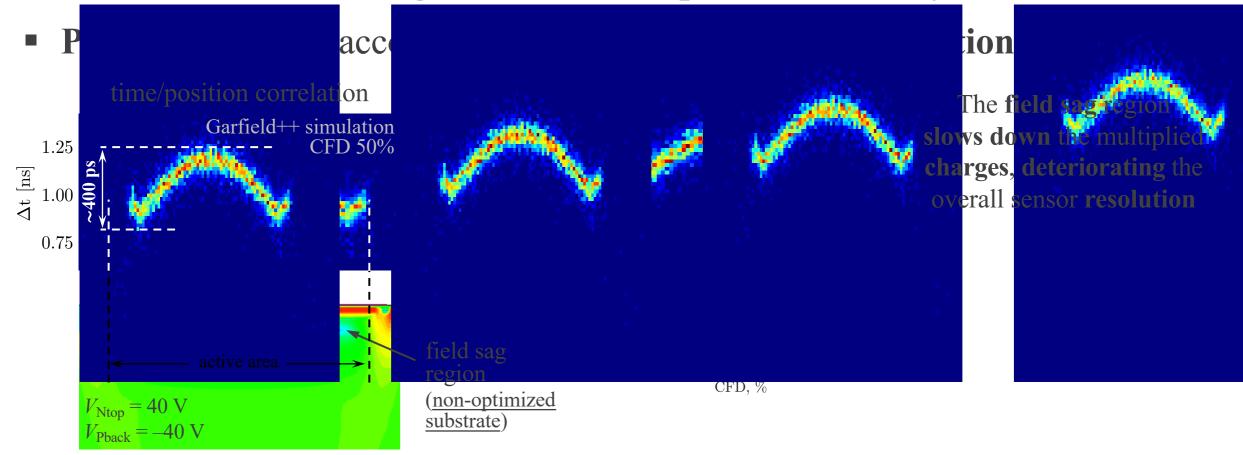
Further numerical investigations on the response uniformity:

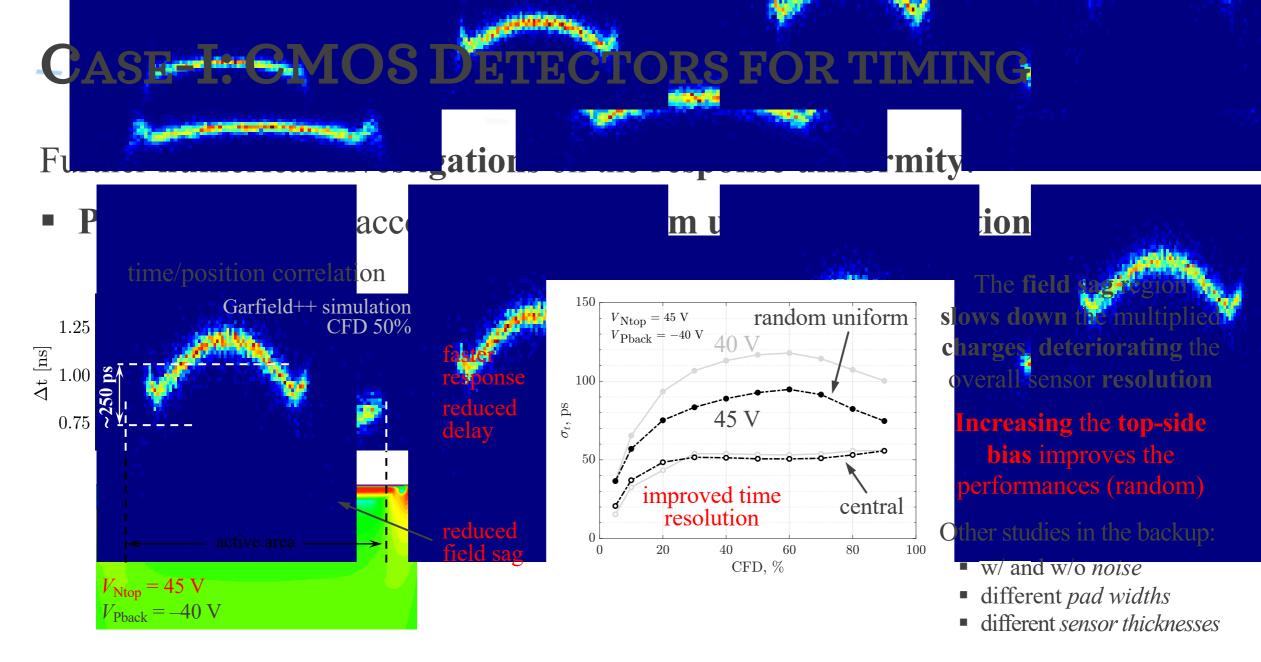
Particle injection according to a random uniform distribution



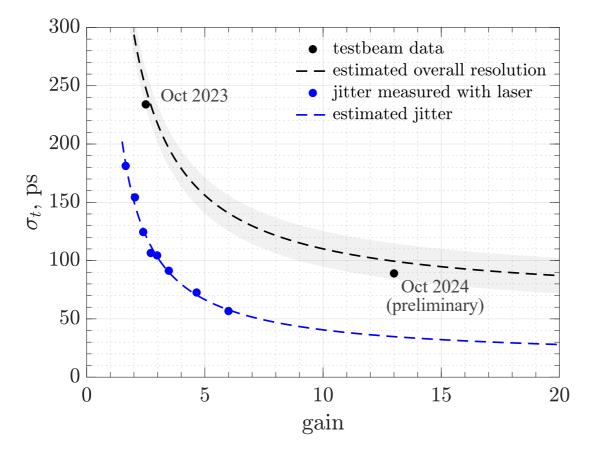
Why the **resolution is high** although **A2** was expected to have **good timing performances**?

#### Further numerical investigations on the response uniformity:



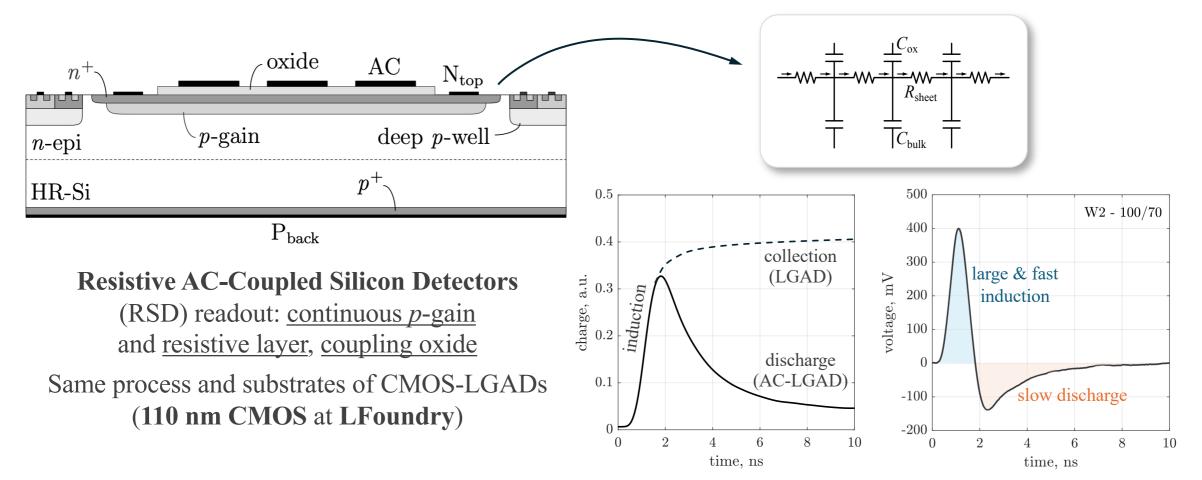


#### Simulated versus measured time resolution in a 50-µm-thick CMOS-LGAD



dashed curves are obtained by fitting **Garfield++ simulations** (*black*) in random uniform distribution configuration and **experimental data** (*blue*)

We are about to integrate the **RSD readout** into the monolithic **CMOS technology** 

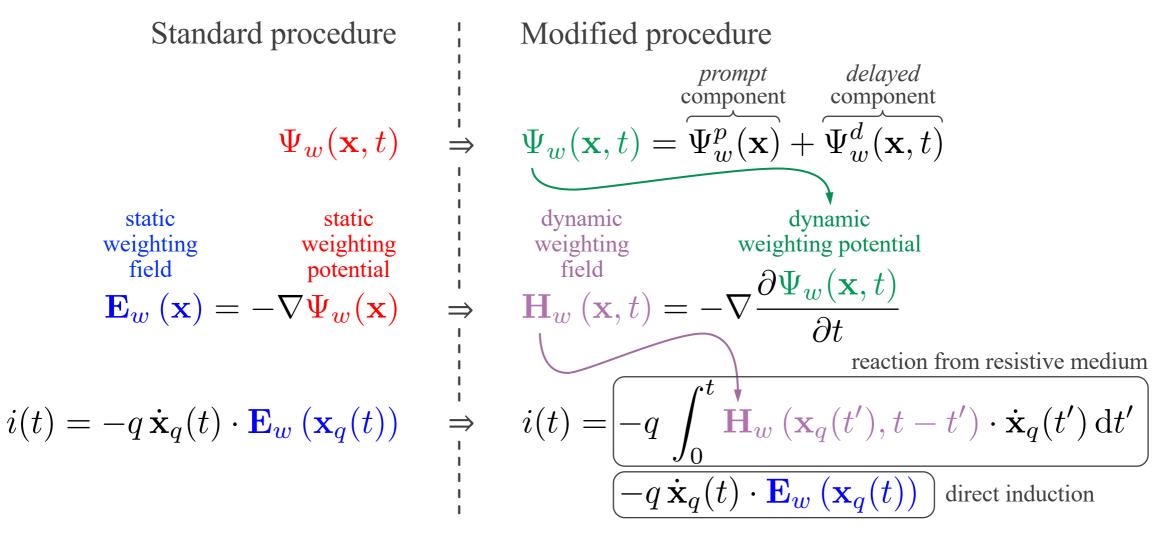


To calculate induced signals in presence of **finite conductivity** elements we need an **extension**<sup>\*</sup> of the **Ramo-Shockley theorem** based on the fact that signal formation depends on both **charge movement** in the **drift medium** and the **time-dependent reaction** of **resistive materials** 

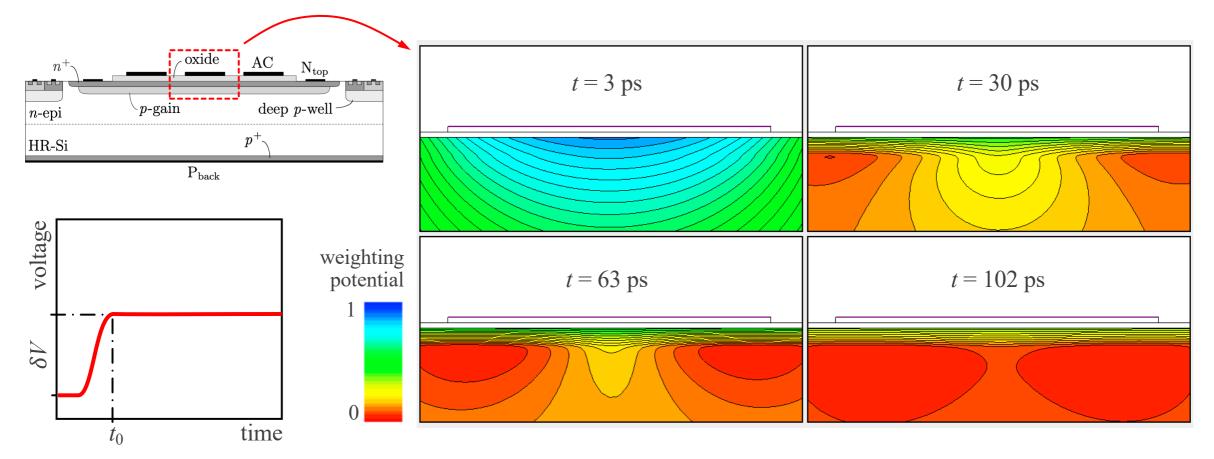
In addition to the usual procedure to extract the potential map, one has to run a **transient simulation** after applying the voltage increment  $\delta V$ , saving the potential at **different points in time**:

$$\begin{cases} V_{\text{bias}} \to \Psi_w^p(\mathbf{x}) \\ \delta V \to \Psi_w^d(\mathbf{x}, t) \end{cases} \Rightarrow \Psi_w(\mathbf{x}, t) = \Psi_w^p(\mathbf{x}) + \Psi_w^d(\mathbf{x}, t), \text{ with } \Psi_w^d(\mathbf{x}, 0) = 0 \end{cases}$$

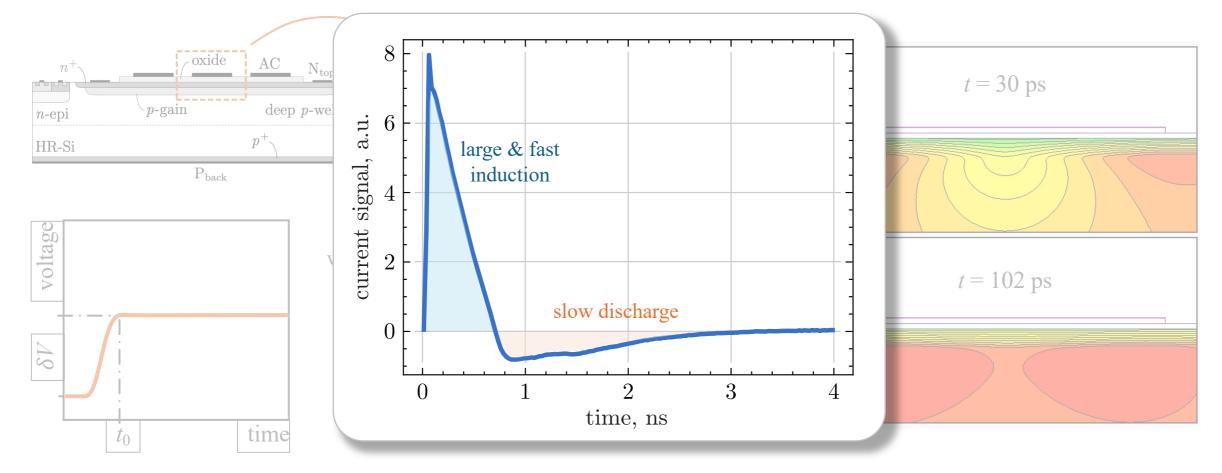
\* see Refs. W. Riegler, NIM A 535 (2004) 287, W. Riegler, NIM A 940 (2019) 453 and D. Janssens, Ph.D. thesis (2024)



A real example of **dynamic weighting potential**  $\Psi_w(\mathbf{x}, t)$  in an **RSD** at work:



#### A real example of **dynamic weighting potential** $\Psi_w(\mathbf{x}, t)$ in an **RSD** at work:





#### FINAL COMMENTS & ACKNOWLEDGMENTS

- Garfield++ is a Monte Carlo toolkit optimized for simulation of ionization, charge transport and current induction in gaseous/semiconductor detectors
- The tool demonstrates very good capabilities in reproducing subtle dynamic effects within the sensor, while significantly reducing computational time with respect to traditional FEM approach
- Garfield++ stands out for some important features like the capability to import all the material parameter libraries and the mesh grid structure from TCAD, as well as to manage the signal induction in structures with resistive elements

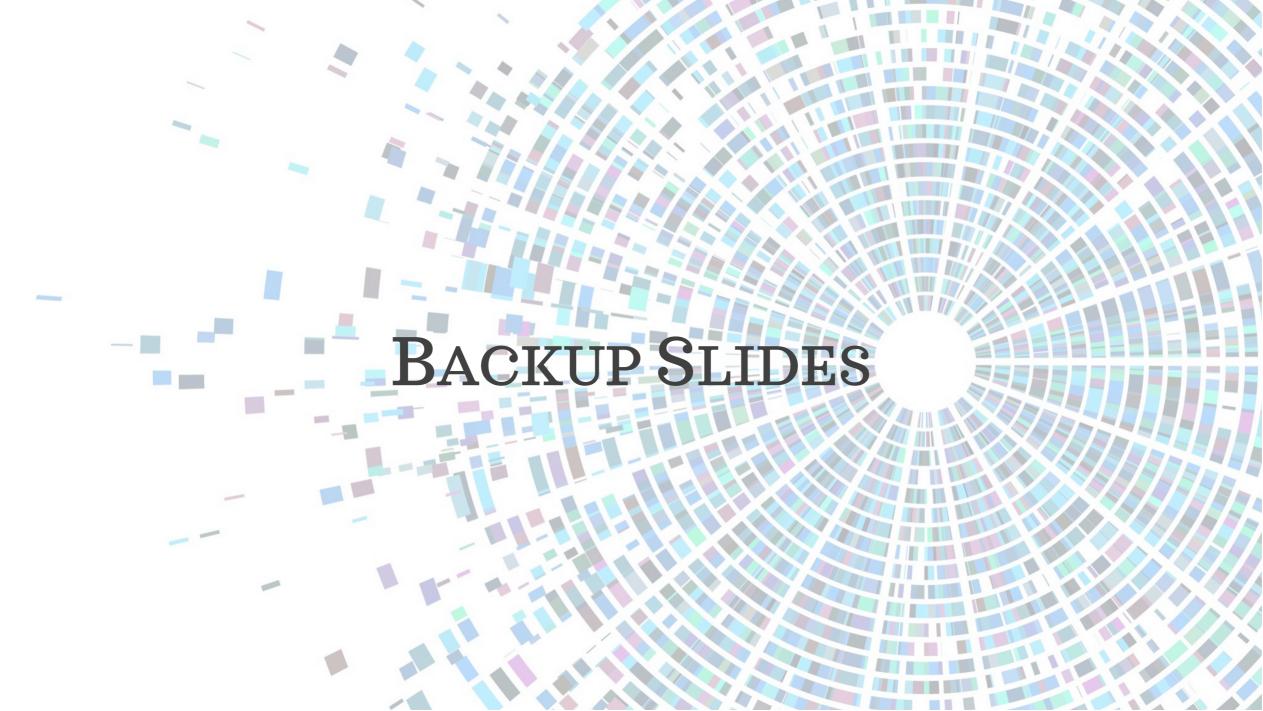
This work would not have been possible without these **funding** and **projects**: **RSD** and **ARCADIA** INFN CSN5 projects, the European Union's Horizon 2020 Research and Innovation Programme, and the **AIDAInnova** programme

A special acknowledgment goes to the following **people** for their help and valuable discussions U. Follo, D. Janssens, H. Schindler, G. Gioachin, G. Andrini, C. Ferrero, S. Durando, L. Pancheri

## THANK YOU FOR YOUR ATTENTION

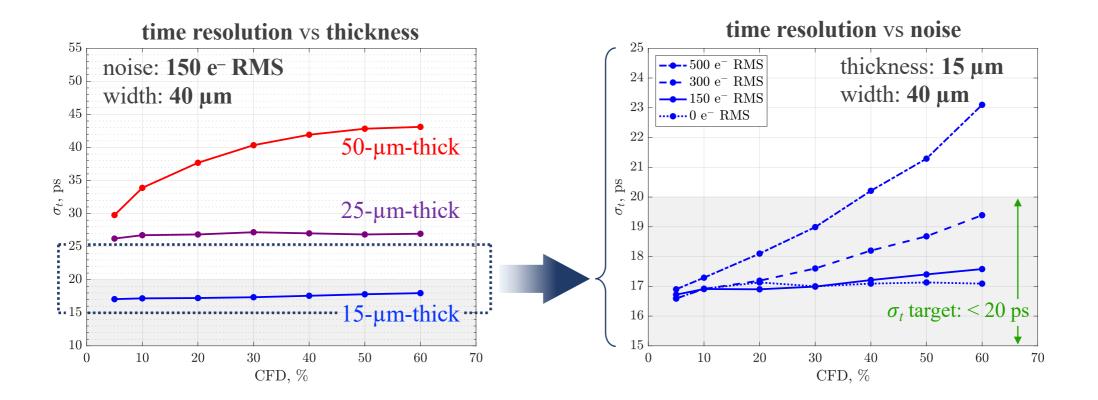
ANY QUESTIONS.?





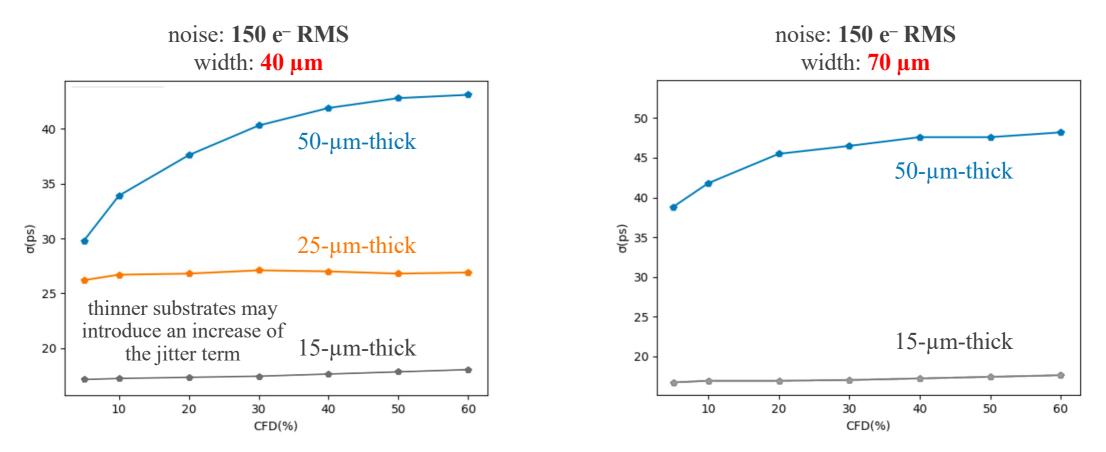
## SIMULATIONS OF TIME RESOLUTION

#### Time resolution vs. sensor thickness and noise



## SIMULATIONS OF TIME RESOLUTION

Time resolution vs. sensor width



## FRONT-END CONTRIBUTION IN GARFIELD++

