HAPSPIDE WP3 Simulation

18/11/2021

WP3 TASKs (description)

- 3.1 Modeling development/validation.
- 3.2 Geant4 simulations.
- 3.3 TCAD simulations: DC, (AC), transient analysis (CCE).
- 3.3 Optimization of a-Si:H devices: layout/geometries, operating conditions.
- 3.4 Radiation tolerance analysis.

WP3 Milestones

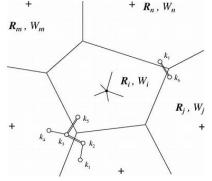
- 3.1 TCAD modeling of a-Si:H M6.
- 3.2 Combined Geant4 + TCAD simulation environment set-up M12.
- 3.3 Comparison between simulations and measurements M18.

WP3 Deliverables / Timescale

- 3.1 a-Si:H TCAD material model with embedded physical models (e.g. carrier mobility) M6.
- 3.2 Report on layout/geometry of dosimetric flux sensor M12.
- 3.3 Report on layout/geometry of single particle sensor M18.
- 3.4 Radiation damage model M36

- WP3 Simulations A twofold framework
 - Dosimetric measurements and single particle detection.
 - Monte-Carlo approach (GEANT4), aiming at accurate description of the material/particle interaction accounting for secondary emission effects, linear energy transfer value variations, effect of packaging materials on tissue equivalency for charge particles and photons, simulation of the expected microdosimetric spectra for different sensitive volume geometries and charge particle beam energy (-> input for TCAD simulations...).
 - Technology CAD (TCAD) approach (Sentaurus Synopsys TCAD), aiming at accurate description of the charge transport within a-Si:H material and device electrical response (DC, transient simulations).
 - The **effect of the read-out mode** (e.g. single-event vs. continuous current measurement) can be self-consistently accounted for within TCAD simulation, exploiting the mixed device-circuit simulation embedded option.

- Innovation: a-Si:H not included in TCAD material library
 - Semiconductor (~ uniform domain macro properties) band gap engineering (defects parametrization).
 - Atomistic scale ? (grain boundary effect, Voronoi domains).



- Charge Transport: mobility models
 - Poor modeling with existing models (e.g. Poole-Frenkel).
 - The (drift) mobility can't be straightforwardly inferred from the current density approximation

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$$J = AT^2 \left[\exp\left(bT\sqrt{\frac{v}{d}}\right) - 1 \right] \iff J = q\mu nE.$$

- Dependence of $\boldsymbol{\mu}$ upon doping, electric field, ...

- Study of LIFETIMES vs. MOBILITY models
 - Interplay in current densities calculations.

- Introduction of innovative materials (from TCAD perspective)
 - Transparent Contact Oxides (TCO) -> transport modeling within oxides.
 - Electrons/Holes selective Contacts -> selective mobility for e, h.

- Simulations vs. measurements @mobility/lifetimes models
 - Current activities/starting point.

