

Use of FLUKA for the analytical calculation of induced radioactivity in the CERN accelerator complex with the JEREMY code

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Outline

Problem
description

FLUKA
approach

JEREMY

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Applications

Conclusions

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2 FLUKA approach

3 JEREMY

4 Implementation

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Problem description

- Beam losses lead to material activation of components of the CERN accelerator complex
- Computation of the nuclide inventory of components of the CERN accelerator complex is necessary
 - Elimination path depends on the nuclide inventory
 - Measurements for all components are too costly (time and money)
- Challenges
 - Large uncertainties in irradiation histories
 - Beam loss maps hardly exist
 - Varying material compositions
 - Isotope production cross sections at high energies

FLUKA approach

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- Full FLUKA simulation with RESNUCLEi scoring
- Pros
 - Correlations are taken into account
 - Detailed description of the geometry and spatial distributions
- Cons
 - Error estimation very difficult / time consuming
 - Production and decay information cannot be easily accessed
 - Long iteration cycle

Spectra + cross section approach

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- The idea is to compute the fluence spectra of the particles in the hadronic cascades via Monte Carlo simulation and fold them with the cross sections offline
- Required inputs
 - Spectra for neutrons, protons, pions and photons
 - Isotope production cross sections
 - Irradiation history
- Output
 - Isotope production rates
 - Specific activity (relative or absolute)
- Extension of PSI approach

Workflow

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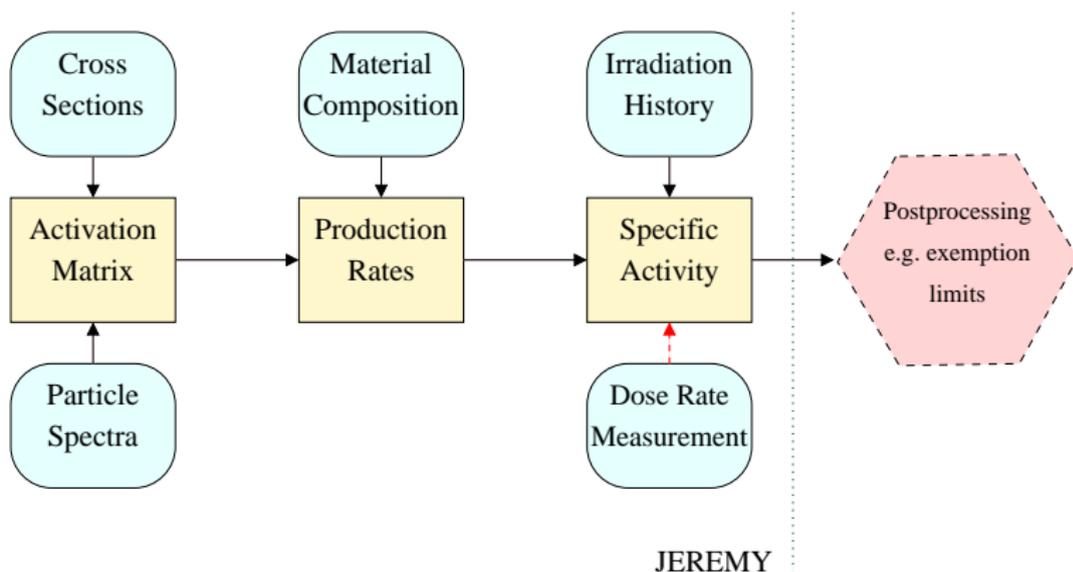
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Inputs obtained with FLUKA

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- Spectra for neutrons, protons, pions and photons
- Isotope production cross sections apart from low energy neutrons (≤ 20 MeV)
 - For isomeres also use FLUKA approach (50% ground state, 50% isomeric state)

Matrix formalism

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- Specific activity

$$A_b = \sum_r \sum_e T_{br} P_{re} m_e$$

- Isotope production rate

$$P_{re} = \frac{N_A}{M_e} \sum_{i=p,n,\gamma,\pi^+,\pi^-} \int \Phi_i(E) \sigma_{i,e,r}(E) dE$$

- Time evolution

$$\begin{aligned} T_{br}(t_{\text{irr}}, t_{\text{cool}}) &= \sum_{c,r \rightarrow b} \int_0^{t_{\text{irr}}} \sum_{m=1}^{j_c} c_{m,c} e^{-\lambda_{m,c} ((t_{\text{cool}} + t_{\text{irr}}) - t_0)} dt_0 \\ &= \sum_{c,r \rightarrow b} \sum_{m=1}^{j_c} \frac{c_{m,c}}{\lambda_{m,c}} \left(e^{-\lambda_{m,c} t_{\text{cool}}} - e^{-\lambda_{m,c} (t_{\text{cool}} + t_{\text{irr}})} \right) \end{aligned}$$

Error estimation

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- Standard Gaussian error propagation for
 - Material composition
 - Irradiation history
 - Cross sections
 - Particle spectra
- Monte Carlo simulation

Standard Gaussian error propagation

- Derivative with respect to the material composition

$$\frac{\partial A_b}{\partial m_e} = \sum_r T_{br} P_{re}$$

- Derivative with respect to t_{cool}

$$\frac{\partial T_{br}}{\partial t_{\text{cool}}} = \sum_{c,r \rightarrow b} \sum_{m=1}^{j_c} -c_{m,c} \left(e^{-\lambda_{m,c} t_{\text{cool}}} - e^{-\lambda_{m,c} (t_{\text{cool}} + t_{\text{irr}})} \right)$$

- Derivative with respect to t_{irr}

$$\frac{\partial T_{br}}{\partial t_{\text{irr}}} = \sum_{c,r \rightarrow b} \sum_{m=1}^{j_c} c_{m,c} e^{-\lambda_{m,c} (t_{\text{cool}} + t_{\text{irr}})}$$

Standard Gaussian error propagation

- Derivative with respect to the l -th bin $\Phi_i(l)$ of the spectrum histogram

$$\frac{\partial P_{re}}{\partial \Phi_i(l)} = \frac{N_A}{M_e} \sigma_{i,e,r}(l) \Delta E_i(l)$$

where $\Delta E_i(l) = \bar{q}_{i,l+1} - \bar{q}_{i,l}$.

- Derivative with respect to the l -th bin $\sigma_{i,e,r}(l)$ of the isotope production cross section histogram for isotope r by shooting particle type i onto element e

$$\frac{\partial P_{re}}{\partial \sigma_{i,e,r}(l)} = \frac{N_A}{M_e} \Phi_i(l) \Delta E_i(l)$$

Standard Gaussian error propagation

- Error on the induced specific activity A_b

$$\begin{aligned}\sigma_{A_b}^2 = & \sum_{t_i, t_j = t_{\text{cool}}, t_{\text{irr}}} \frac{\partial A_b}{\partial t_i} \sigma_{t_i, t_j} \frac{\partial A_b}{\partial t_j} + \sum_{m_{e_i}, m_{e_j}} \frac{\partial A_b}{\partial m_{e_i}} \sigma_{m_{e_i}, m_{e_j}} \frac{\partial A_b}{\partial m_{e_j}} + \\ & \sum_{i=p, n, \gamma, \pi^+, \pi^-} \sum_{\Phi_i(l), \Phi_i(j)} \frac{\partial A_b}{\partial \Phi_i(l)} \sigma_{\Phi_i(l), \Phi_i(j)} \frac{\partial A_b}{\partial \Phi_i(j)} + \\ & \sum_r \sum_e \sum_{i=p, n, \gamma, \pi^+, \pi^-} \sum_{\sigma_{i, e, r}(l), \sigma_{i, e, r}(j)} \frac{\partial A_b}{\partial \sigma_{i, e, r}(l)} \sigma_{\sigma_{i, e, r}(l), \sigma_{i, e, r}(j)} \frac{\partial A_b}{\partial \sigma_{i, e, r}(j)}\end{aligned}$$

- Full covariance formalism possible
 - Correlations between spectrum bins computed with FLUKA would be nice
- Impact of individual quantities (e.g. the cross section for a given reaction) can be assessed
- Error propagation for derived quantities

Monte Carlo simulation

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- Assign PDFs to relevant parameters according to their uncertainty and obtain distributions for the specific activity
- in principle arbitrary PDFs
- all common PDFs available, e.g. gaussian, exponential
- Very fast approach (2000 events per minute)
- Distributions of derived quantities can be easily computed

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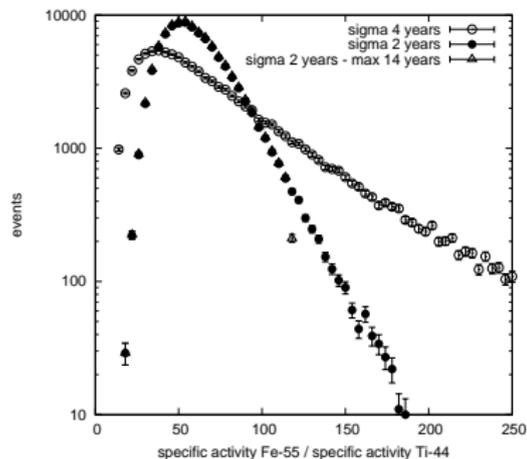
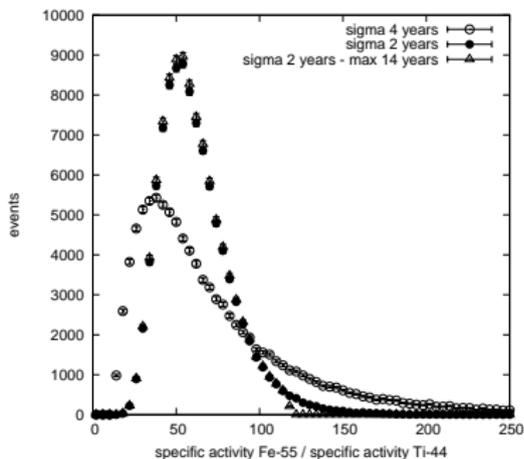
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Implementation of JEREMY

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- Coded in Python with `numpy` package
 - Selected core routines coded optionally in C for speed-up
- Documented in `pydoc/html` format
- Object oriented
 - in combination with Python → powerful scripting interface, e.g. for postprocessing (exemption limits) or Monte Carlo simulation
- Isotope production cross sections
 - Low energy neutrons from JEFF 3.1.1 library (including isomere branching ratios)
 - High energy particles from FLUKA models
- Full time evolution with decay data from JEFF 3.1.1 library
- GUI (planned)

Applications of JEREMY

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- Characterization and elimination of first components of the CERN SPS complex towards France this year
- Computation of the nuclide inventory for components that will be eliminated towards Switzerland for the capacity planning for the Swiss Final Repository
- Radiological classification of parts of LHC detectors
- Hadron therapy centers

Conclusions

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- Improved uncertainty estimation for induced activity
- Full access to production and decay information
- FLUKA simulations heavily needed for spectra and cross sections