

# Using Parallel Monte Carlo to Validate Radiotherapy Calculations

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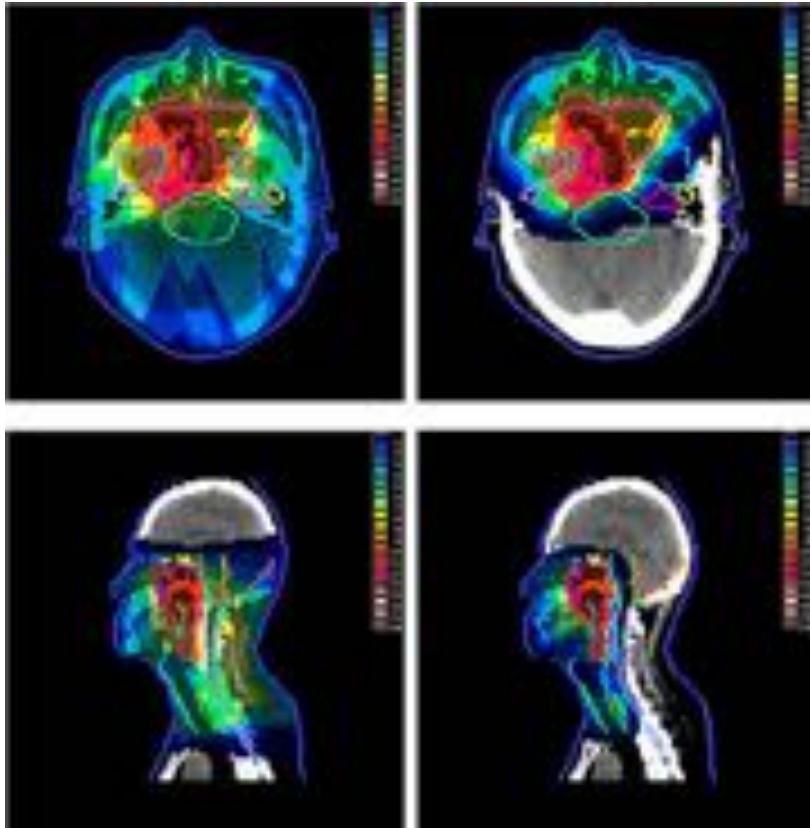
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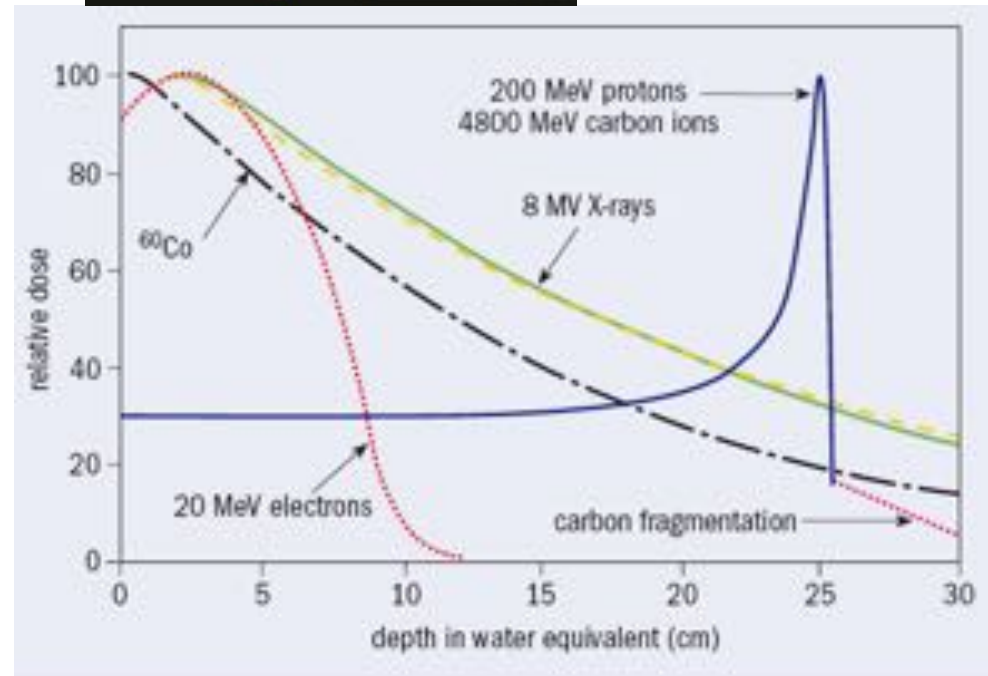
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## Protons vs X-Rays



IMRT  
Intensity  
Modulated  
Radiotherapy



$$-\frac{dE}{dx} = \frac{4\pi}{m_e c^2} \cdot \frac{n z^2}{\beta^2} \cdot \left( \frac{e^2}{4\pi\epsilon_0} \right)^2 \cdot \left[ \ln \left( \frac{2m_e c^2 \beta^2}{I \cdot (1 - \beta^2)} \right) - \beta^2 \right]$$

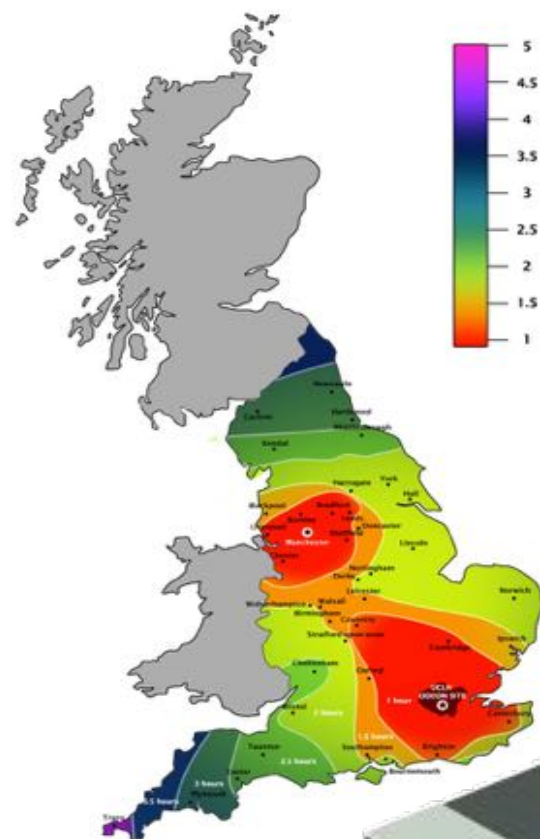


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MANCHESTER  
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The University of Manchester

# These centres are big...

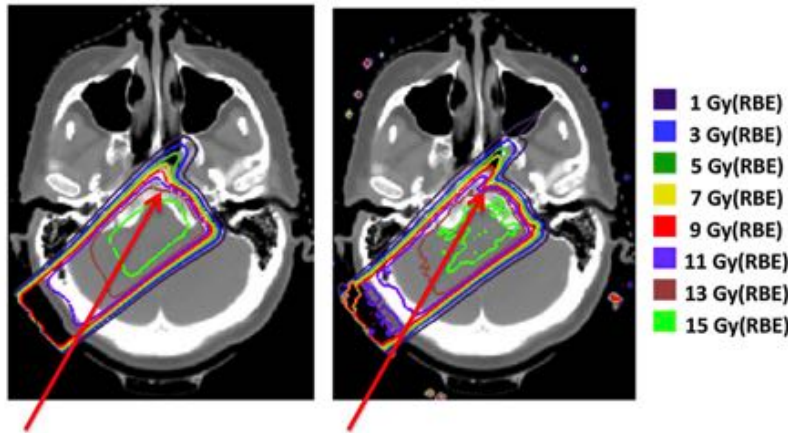


£125M each  
750 patients per year  
(~30 treatments per patient)  
£25k per patient  
18 hours running per day  
Running from 2018 onwards

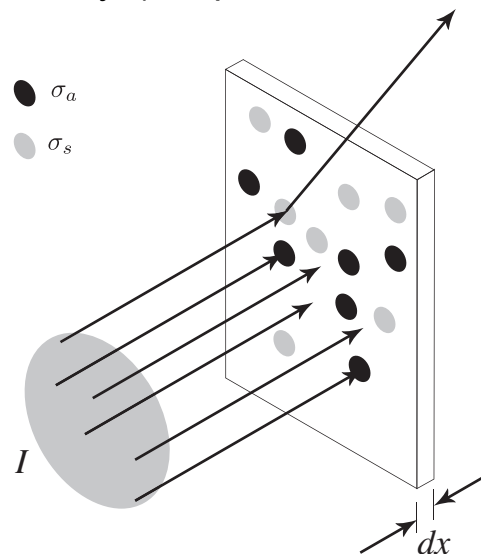
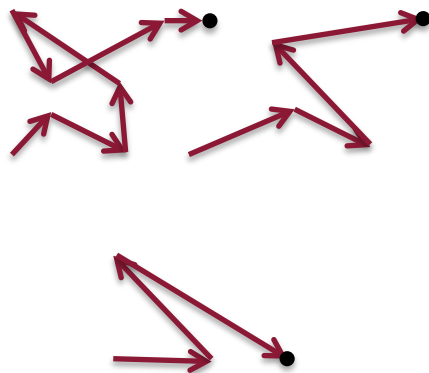


UCLH Centre

# The Role of Monte Carlo in Proton Therapy



Pencil beam algorithm (left) vs Monte Carlo (right); arrow indicates a range error due to MCS in heterogenous boundary. (Adapted from Paganetti (2008))



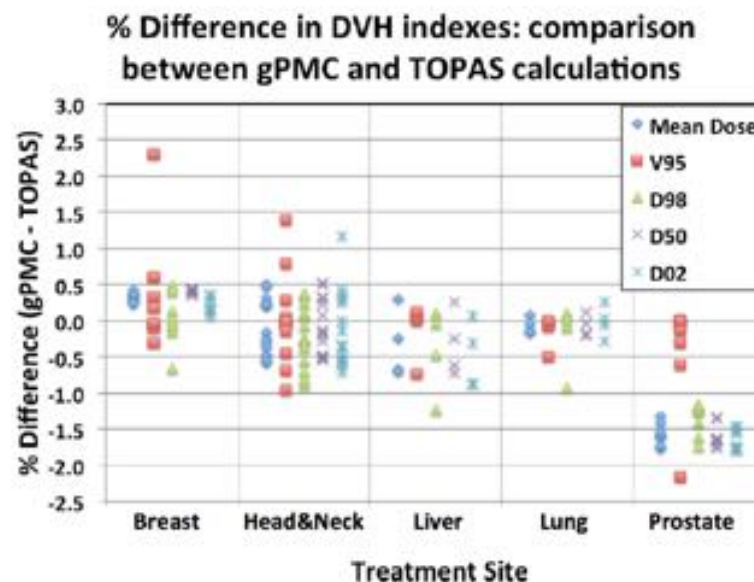
- Algorithms used in planning tools can have drawbacks.
- Monte Carlo is too slow to be used directly for planning (maybe not, see later...)
- Use Monte Carlo to validate plans before delivery
- Must fit planning workflow:
  - Import patient CT file as a geometry of voxels.
    - Usually a few million, more in head and neck
    - c. 40 hours on single CPU
  - Track protons through the voxels recording where the energy goes.
    - Must simulate enough protons to get good uncertainty (< 1%).
    - Must be confident in the physics implementation.
    - Needs to fit within a clinical workflow – minutes not hours.





## gPMC – Simplified Physics using GPU

- Image grid 1 x 1 x 1.5 mm
- Dose grid 2 x 2 x 2.5 mm
- Simulation rate
  - gPMC: ~2.6s/MP
  - TOPAS: 4h/MP
- Simulation time
  - 50-500 MP (histories)
  - gPMC: ~130s to 1300s
- Random uncertainties (depends on site):
  - gPMC: 0.5-2.4%
  - TOPAS: 1-2%
- Dose difference gPMC/TOPAS:
  - Around 1% on D98, D50, D02
- This problem is in the simulation of nuclear interactions, exacerbated by the higher energies required in prostate treatment

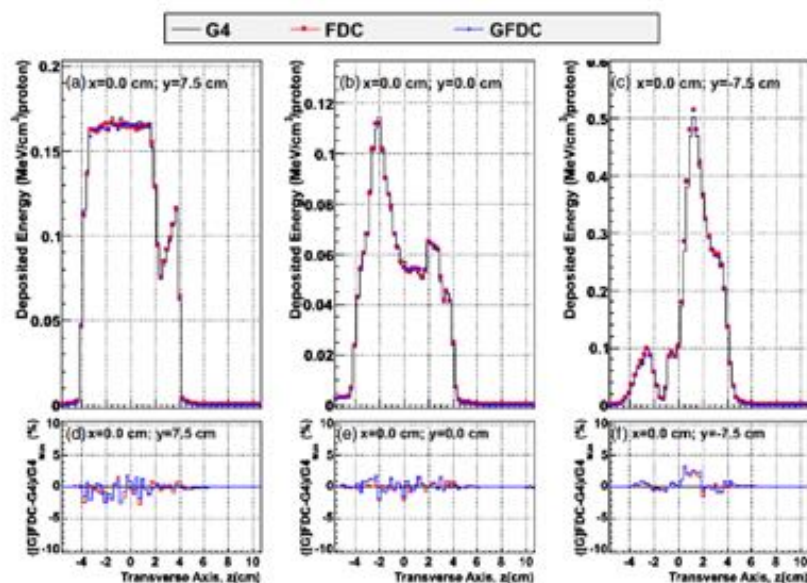


Giantsoudi *et. al.* PMB60, 2257 (2015)

- *gPMC can already calculate good accuracy in a few minutes!*



## Track Repeating



Dose distribution projected into each plane. Differences between Geant4 and the GPU track repeating implementation are around 2 %

- Create a database with 1 million proton histories.
- Stores step length, angle relative to previous step, energy lost and energy deposited for a 251 MeV beam in water and 41 other biological materials.
- This data is then scaled for different energies.
- Validated against Geant4 – accurate to ~2%
- 5.4s/MP (on older hardware)



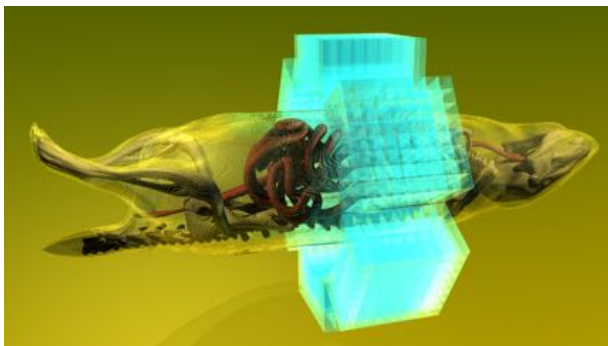
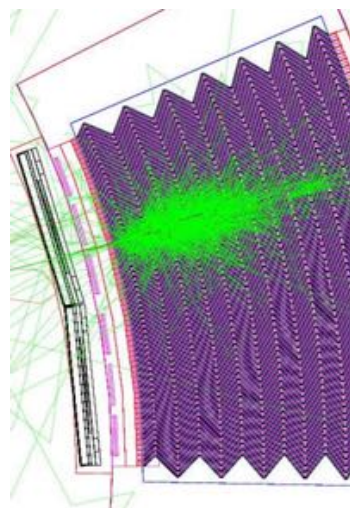
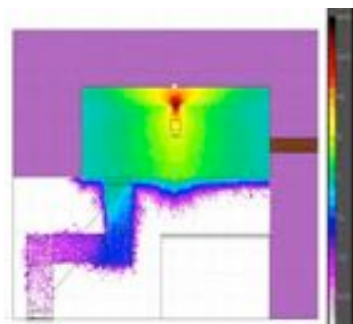
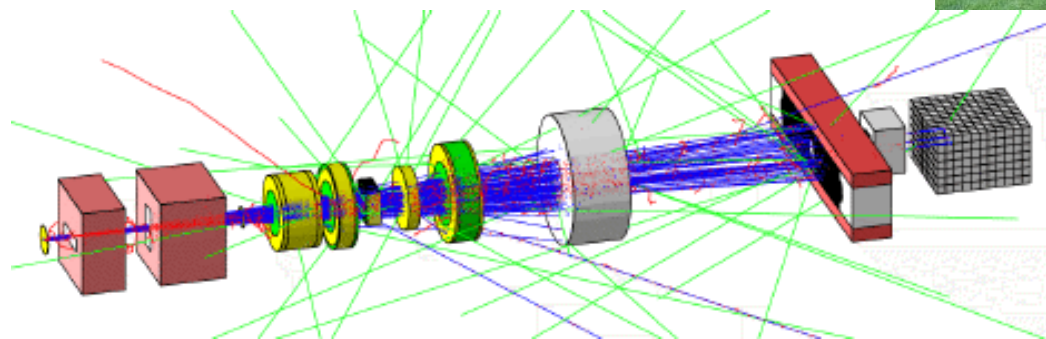
## What about Intel?

- XeonPhi MIC – Many Integrated Core
- 61 cores/244 threads (today)
- Only 16GB on-card RAM (today)
- Challenging target of <67MB/thread.
- **BUT:** x86 architecture
  - Same source code
  - Same compilers
  - Same development cycle
  - Less debugging!
  - Keep all physics!



Model	3120A	7120P
# Cores (threads)	57 (228)	61 (224)
RAM	6 GB	16 GB
Clock Speed	1.100 GHz	1.238 GHz
Cost	\$1695-\$1960	\$4129

*Used Manchester Research IT test system (thanks!)*



Well known in HEP, well validated  
(see e.g. (Yarba, 2012))

Lots of front ends for Medical  
Physics (TOPAS, GATE, GAMOS  
etc)

- Validation in medical applications  
is good (e.g. (Testa *et. al.* 2013))
- Recent addition of multithreading  
opens the possibility of running  
on Xeon Phi
  - Per process memory required is  
~100s MB
  - Per thread can be ~10s MB
- Useful to have a common code  
base on all systems
- We used **stock** Geant4MTv10

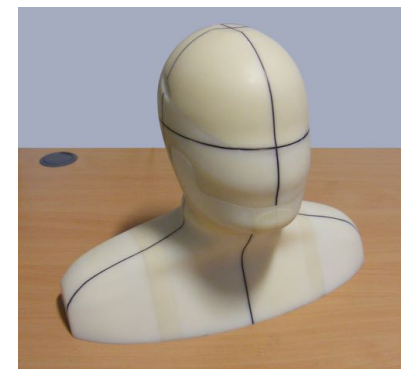




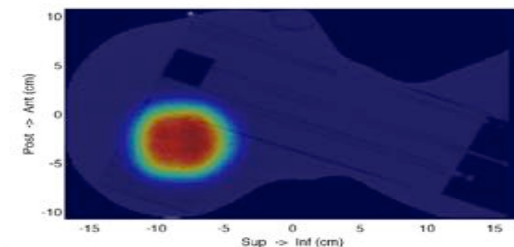
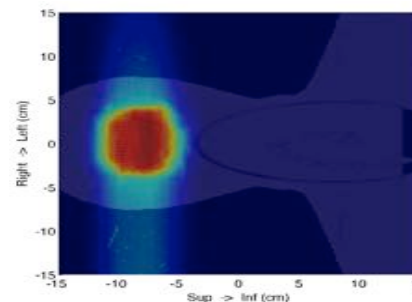
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## Benchmark Calculation

- Used a radiotherapy phantom (Aitkenhead *et. al.* 2013) to avoid data protection issues
- CT scan image converted to materials 16M voxels
- Image grid 1 x 1 x 2 mm
- Dose grid 2 x 2 x 2 mm
- Simulation time
  - 50-500 MP (histories)
  - gPMC: ~130s to 1300s
- Two beam angles, each with c.1500 spots;  
~3000 GPS sources
- Simulation of 10MP:
  - ~1% uncertainty in high dose region



- The Geant4 General Particle Source is the obvious choice for simulating spots:
  - Any energy distribution.
  - Any source shape/size.
  - Any particle.
  - Weighted sampling.
- However, not previously optimized for multithreading – lots of memory required.

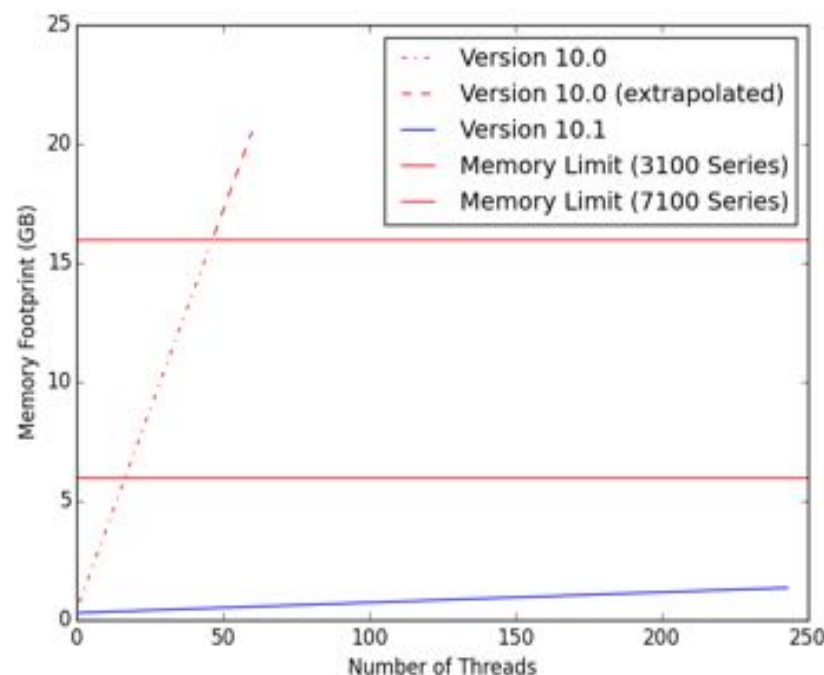


*The dose distribution:  
50 cm<sup>3</sup> sphere roughly 7 cm below surface of head*



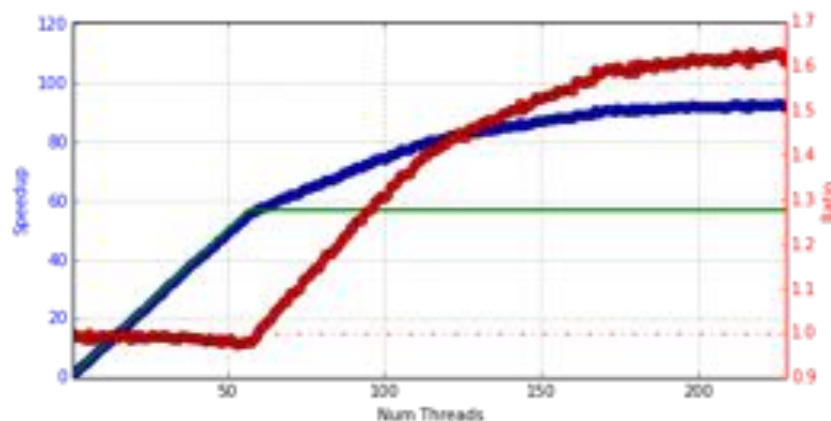
## The Geant4 General Particle Source

- In active scanned proton therapy, need to model many beamlike sources.
  - Simulated by Gaussian distributions in space and energy.
  - The location and shape of the Gaussians define the treatment plan.
- Share data between threads:
  - Things that don't change during a run
  - Source positions, energies, intensities etc.
- Use an approximation to neutron cross sections – negligible impact on accuracy (see Asai *et. al.* 2014)





## Calculation Speed



Speedup as a function of number of threads. Ratio shows the speedup over the number of physical cores – it should be 1 until hyper-threading starts.

- Now the simulation can run, how fast does it go?
  - One 7120P card – 1h 54m for 10MP
  - Two 7120P cards – 1h 1m
  - Best performance – 366 s/MP
  - (cf. 2.6s/MP for gMPC)
  - 2 x XeonPhi is \$13,500
- Speed Comparisons:
  - AMD Opteron machine (48 threads, 2.6 GHz) gives similar speed at \$6,000
  - Single card gPMC about 100x faster

*Bottom Line: Lose about 100x speed cf. GPU, but keeps all physics/accuracy*

## Cloud Computing

- Been done before using Microsoft Azure cloud
- Previous work has looked at cost
  - We also looked at speed
- We used Google Cloud's Google Compute Engine (GCE)
  - Competitive pricing
  - Powerful machines
  - Extremely fast network
- MC codes are almost ideal problem for cloud computing:
  - Independent calculations combined at end

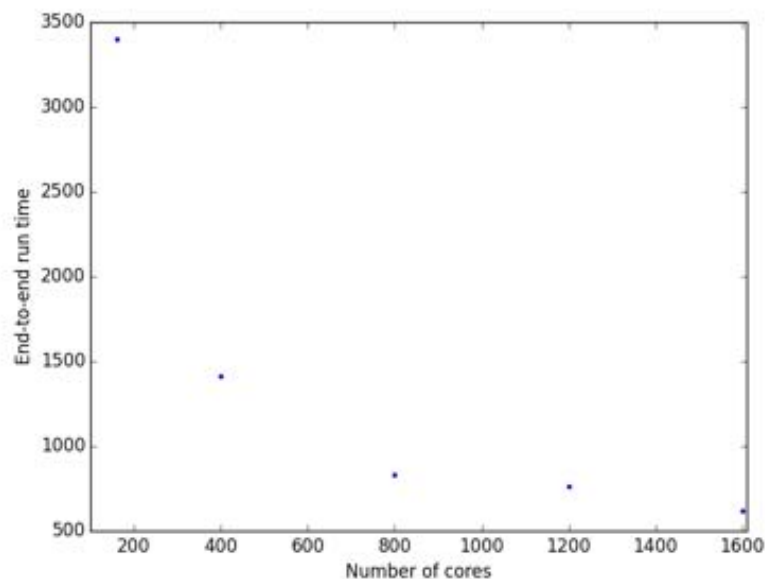






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## Cloud Computing



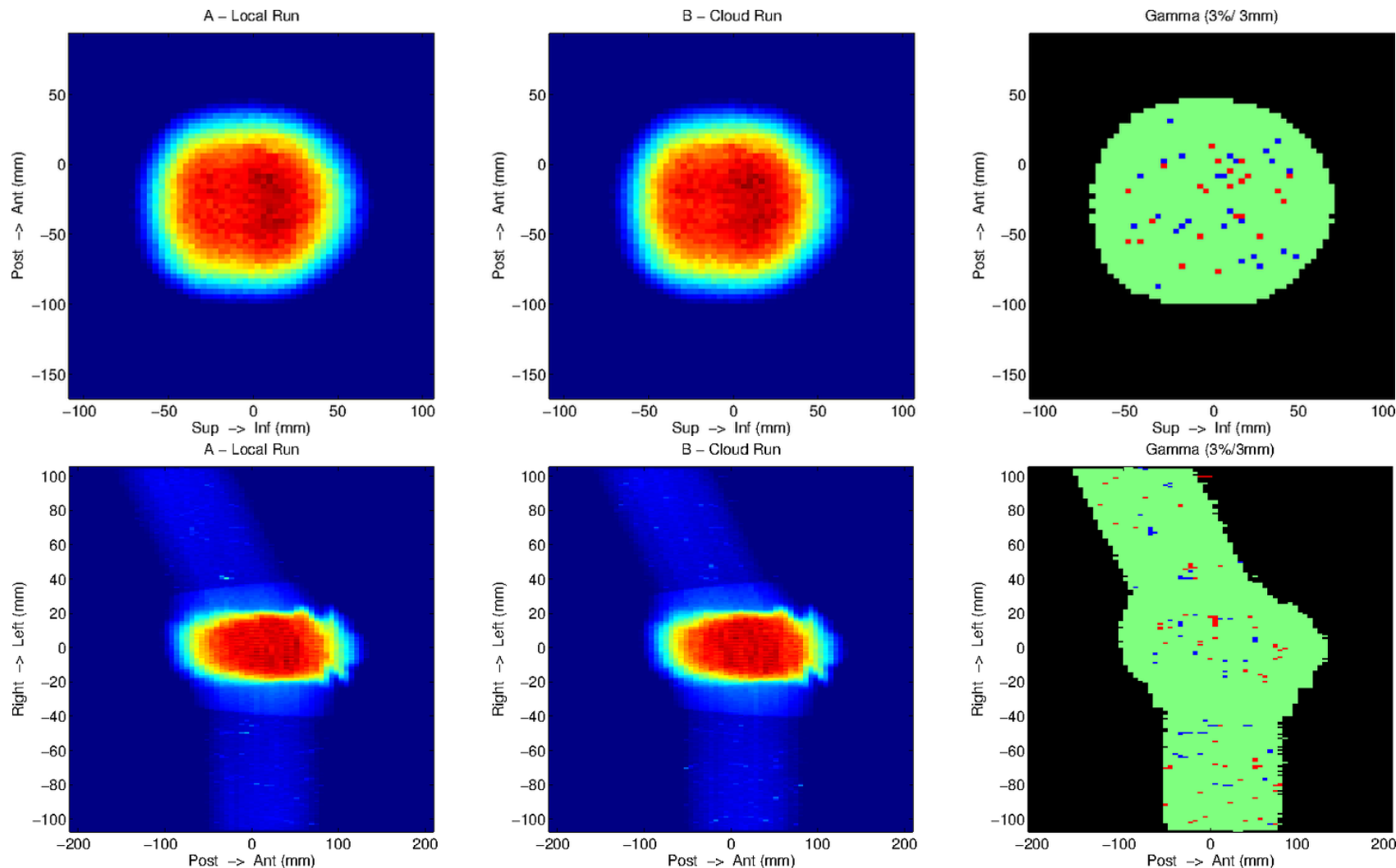
Scaling with number of cloud cores used. Includes calculation of uncertainty and collation and combination of data. Download of the result adds roughly 10 seconds.

- A simple plan requires  $10^7$  proton primaries
  - 36 hours on a desktop
  - 1 hour on 2x Xeon Phi
  - 10 minutes on GCE
  - Roughly on par with gPMC
- Rent a small supercomputer for 10 minutes
  - 1600 cores
- Scales well, and very cost competitive
  - $\ll \$10$  per validation
  - Price dropping fast

More details: <https://epubs.stfc.ac.uk/work/12273478>



## Results – Dose Comparison

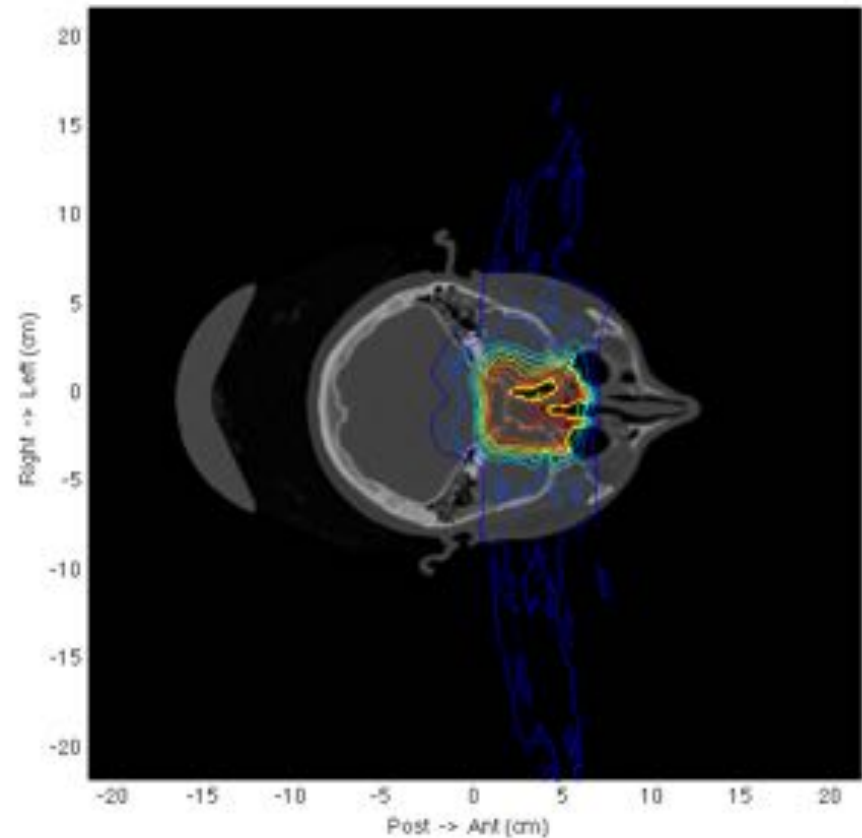


*Gamma analysis shows identical calculation; limited by expected statistical fluctuations*



## Cloud Computing in Radiotherapy

- Issues of patient data control
  - Probably surmountable
  - Needs interpretation/change in NHS data policy
- Cloud will get faster + cheaper
- Especially good fit for hospitals
  - No specialist local skills
  - No specialist local infrastructure
  - Continuous upgrading
  - Highly scalable
- A number of groups/ companies are now looking at this



- Further work on GCE
  - Keep optimising the launch code to minimise overheads
  - Develop MPI version of the code?
  - Completely general: applicable to any other use of GEANT4
- Develop tailored proton therapy application optimised for Xeon Phi
  - Should be faster, but much less general



## Summary

- Geant4 version 10.1 onwards will be able to run proton therapy treatment plans on Xeon Phi
  - Largely due to our memory reductions in the code, especially the GPS.
- A computer with 2 Xeon Phi cards is roughly equal to a 'normal' cluster machine
  - The cluster machine was not latest generation – expect better performance on Intel E5s
  - Expect the next Xeon Phi generation to be about 3x faster
  - This is with essentially no optimisation other than in memory use
- Algorithm changes in the code, and variance reduction techniques may have big benefits
  - Variance reduction is under-studied
- Computation in the cloud is likely to be a big growth area for hospitals
  - Cheaper to let Google do the hosting and maintenance
  - **As fast as gPMC, but with full (well validated) physics**

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USolids library - <http://aidasoft.web.cern.ch/USolids>