



GRAPHCORE IPU INTRO

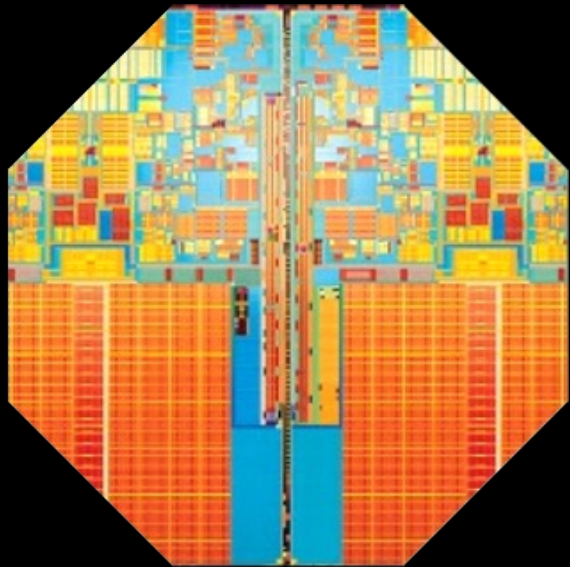
GRAPHCORE

Dr Alex Titterton
Solutions Architect



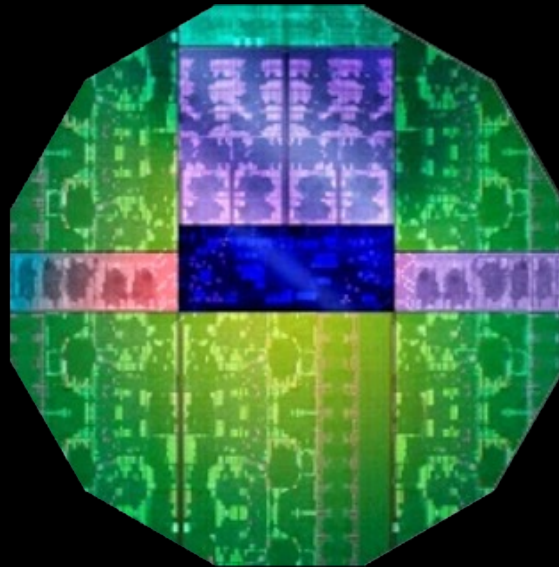
Neural network visualization from [POPLAR™](#)

INTELLIGENCE PROCESSING UNIT DESIGNED FOR AI



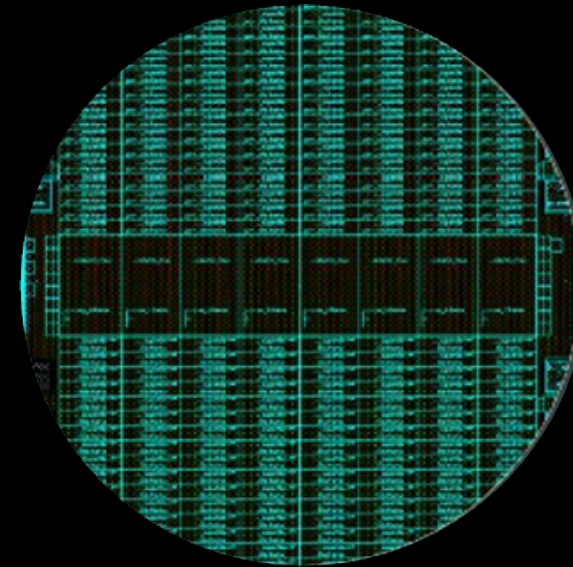
CPU

Scalar



GPU/TPU

Vector



IPU

Graph

THE INTELLIGENCE PROCESSING UNIT (IPU) WHAT MAKES IT DIFFERENT?

CPU

GPU

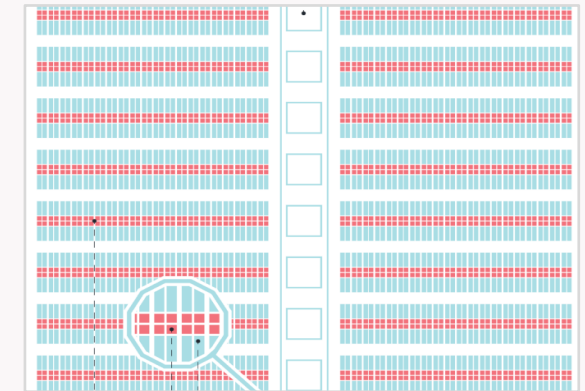
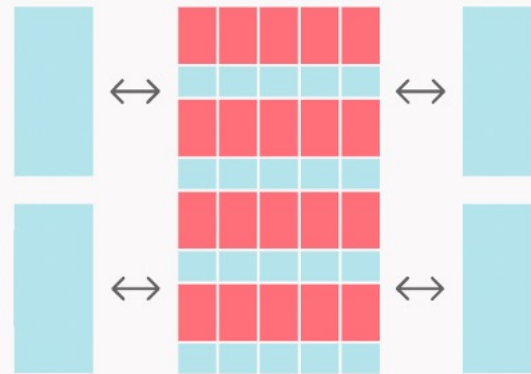
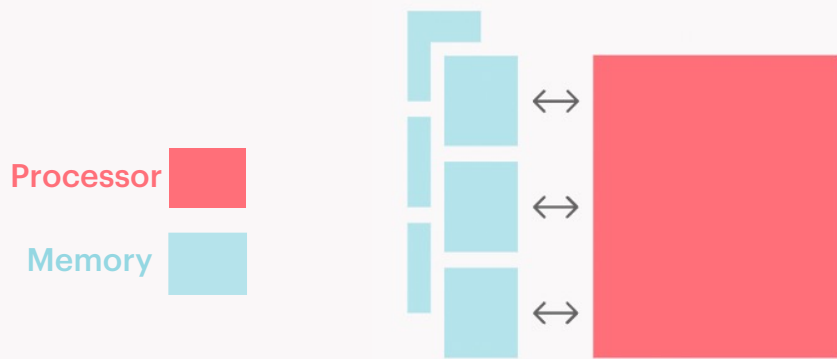
IPU

Parallelism

Designed for scalar processing

SIMD/SIMT architecture. Designed for large blocks of dense contiguous data

Massively parallel MIMD architecture. High performance/efficiency for future ML trends



Memory Bandwidth

Off-chip memory

Model and Data spread across off-chip and small on-chip cache and shared memory

Main Model & Data in tightly coupled large locally distributed SRAM

(2TB/s for A100 HBM)

(~65 TB/s for Bow IPU)

INTRODUCING THE BOW IPU

WORLD'S FIRST 3D WAFER-ON-WAFER PROCESSOR



3D silicon wafer stacked processor

350 TeraFLOPS AI compute

Optimized silicon power delivery

0.9 GigaByte In-Processor-Memory @ **65TB/s**

1,472 independent processor cores

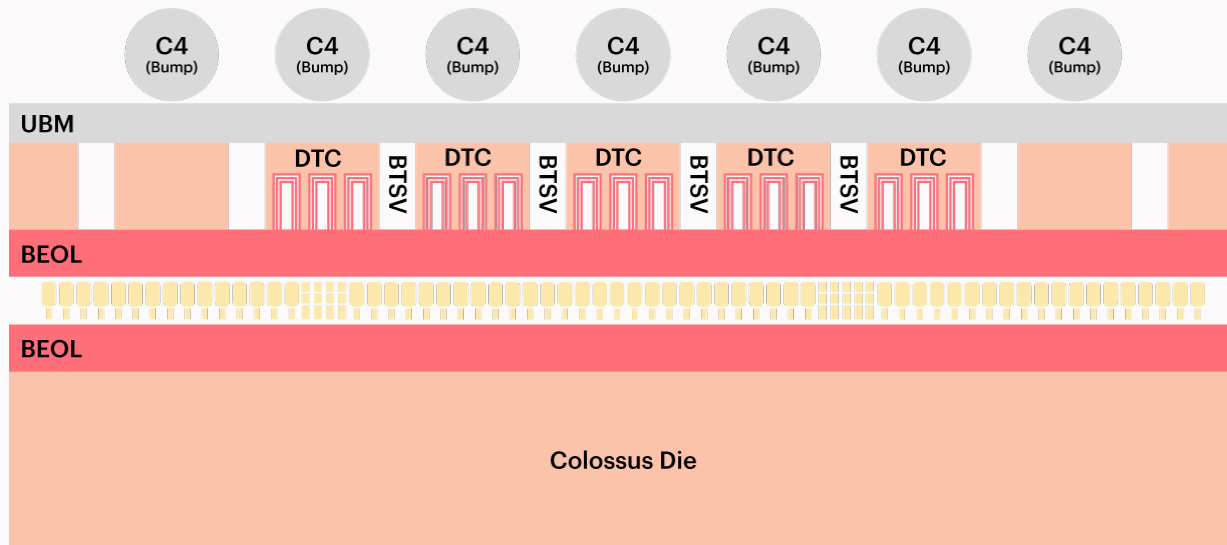
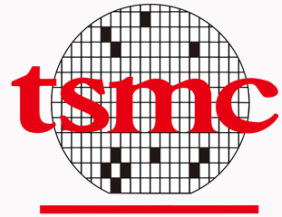
8,832 independent parallel programs

10x IPU-Links™ delivering 320GB/s

BOW IPU: 3D WAFER-ON-WAFER PROCESSOR

GRAPHCORE

+



Advanced silicon wafer stacking technology co-developed between Graphcore and TSMC

World's first commercial deployment using TSMC SoIC-WoW™ technology in Bow IPU

Enabling technology for closely coupled power delivery die to maximize application performance



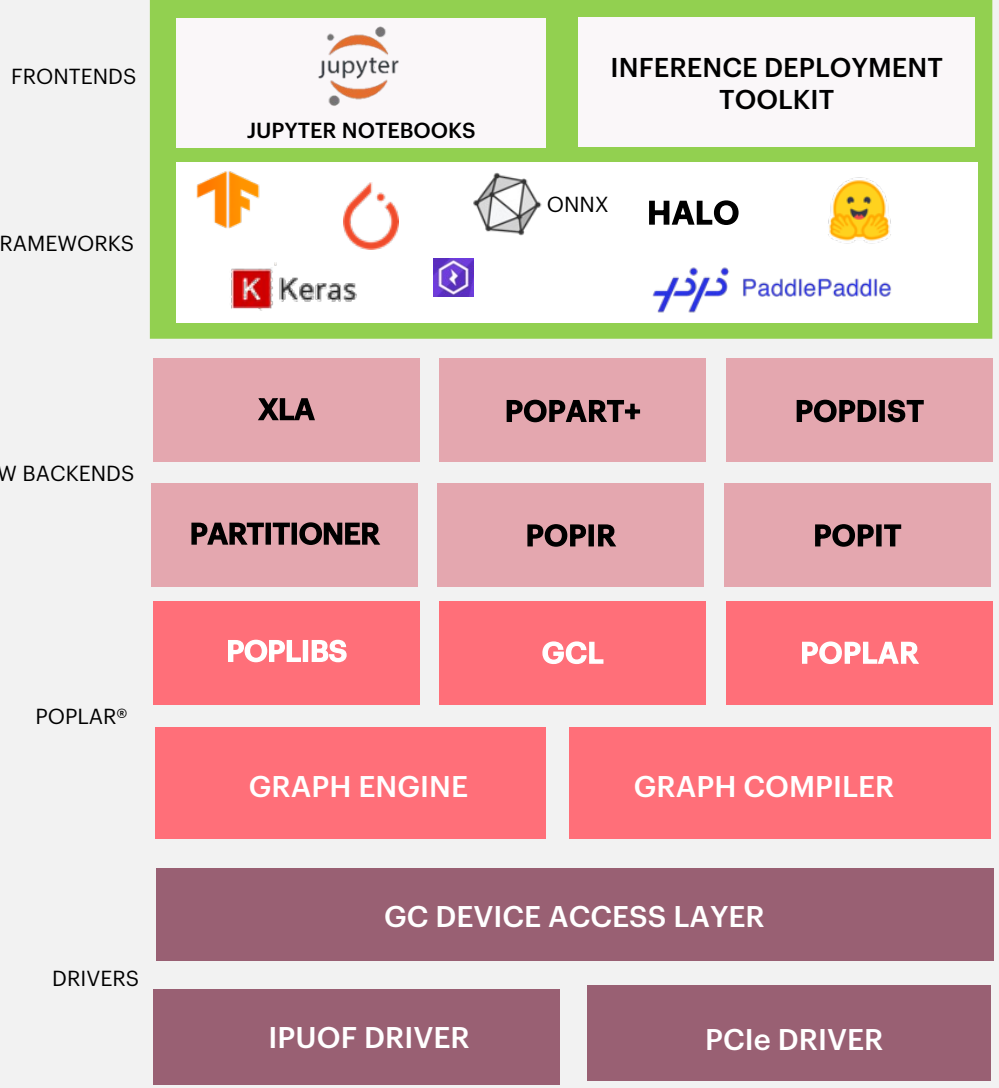
GRAPHCORE SOFTWARE

- NLP/TRANSFORMERS
- IMAGE CLASSIFICATION/CNNS
- OBJECT DETECTION
- LARGE MODELS
- MLPERF
- CONDITIONAL SPARSITY
- GNNS

ML APPLICATIONS

- TUTORIALS
- CODE EXAMPLES
- DOCUMENTATION
- VIDEOS
- NATIVE IPU CODERS PROGRAM
- APPS PORTFOLIO

DEVELOPER ECOSYSTEM



POPULAR® SDK

- GRAPH ANALYZER
- SYSTEM ANALYZER
- DEBUGGER
- DEVELOPMENT ENVIRONMENT

POPVISION TOOLS

- V-IPU
- SYSTEM MONITORING
- PROMETHEUS
- GRAFANA

JOB DEPLOYMENT

- K8S
- SLURM
- SLURM

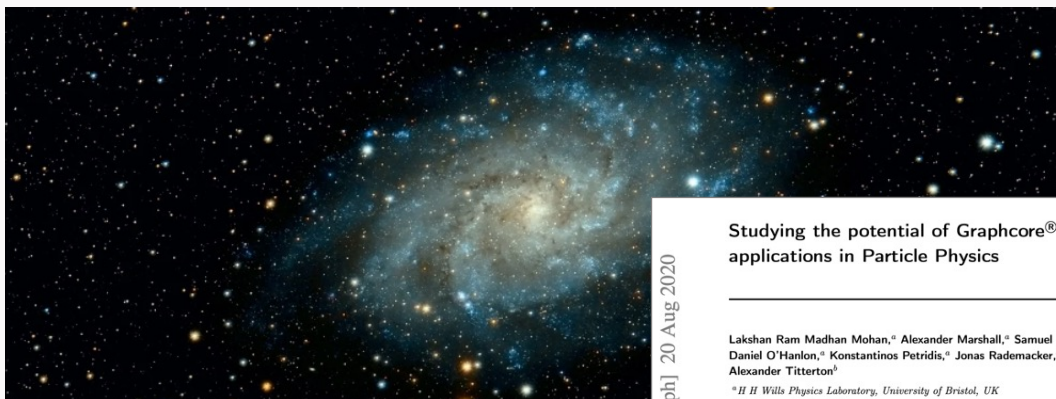
SYSTEM SOFTWARE



GRAPHCORE RESEARCH COLLABORATIONS (SMÖRGÅSBORD)



IPUs in Research



UNIVERSITY OF BRISTOL TACKLES CHALLENGES IN PARTICLE PHYSICS WITH GRAPHCORE'S IPU

arXiv:2008.09210v1 [physics.comp-ph] 20 Aug 2020

Studying the potential of Graphcore® IPU for applications in Particle Physics

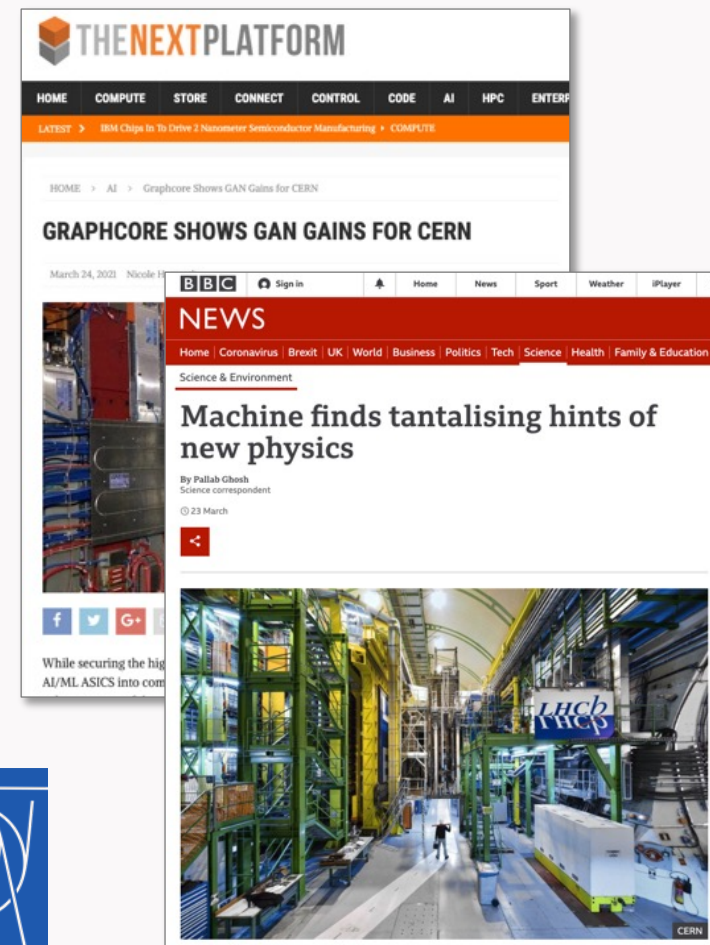
Lakshan Ram Madhan Mohan,^a Alexander Marshall,^a Samuel Maddrell-Mander,^{a,b} Daniel O'Hanlon,^a Konstantinos Petridis,^a Jonas Rademacker,^a Victoria Rege,^b and Alexander Titterton^b

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ABSTRACT: This paper presents the first study of Graphcore's Intelligence Processing Unit (IPU) in the context of particle physics applications. The IPU is a new type of processor optimised for machine learning. Comparisons are made for neural-network-based event simulation, multiple-scattering correction, and flavour tagging, implemented on IPUs, GPUs and CPUs, using a variety of neural network architectures and hyperparameters. Additionally, a Kálmán filter for track reconstruction is implemented on IPUs and GPUs. The results indicate that IPUs hold considerable promise in addressing the rapidly increasing compute needs in particle physics.



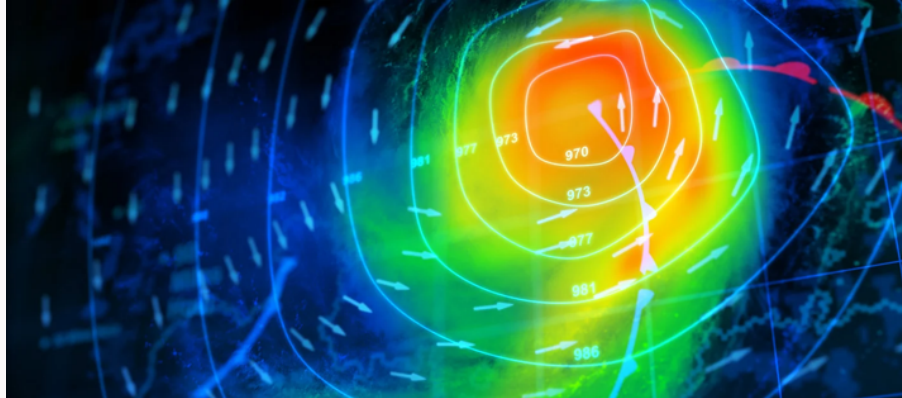
<https://www.graphcore.ai/resources/research-papers>

Using AI to accelerate HPC Scientific Applications

Mar 09, 2022 \ AI, HPC

AI FOR SIMULATION: HOW GRAPHCORE IS HELPING TRANSFORM TRADITIONAL HPC

Written By:
Alex Titterton



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Get Updates

For many years High Performance Computing (HPC) techniques have been used to solve the world's most complex scientific problems across a wide range of applications, from modelling Higgs boson decay at the Large Hadron Collider to using Monte-Carlo simulation to predicting whether the weather will improve.

However, due to the immense complexity of the calculations involved in many of these applications, researchers are often waiting a long time for simulation results to arrive. Speeding up these workflows by simply running the same programs on more powerful hardware can be very expensive, with a large cost often giving only a modest improvement in performance.

Clearly, a new approach is required to efficiently speed up these workloads, and many researchers are turning to surrogate machine learning models.

A surrogate model is a machine learning model intended to imitate part of a traditional HPC workflow,

For more information, see our technical blog post:

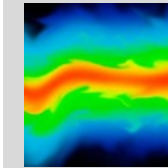
<https://www.graphcore.ai/posts/ai-for-simulation-how-graphcore-is-helping-transform-traditional-hpc>



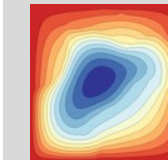
Relevant Application Areas



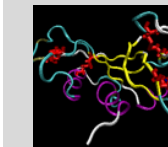
High Energy Physics



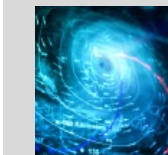
Computational Fluid Dynamics



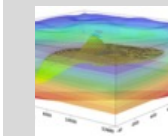
Partial Differential Equations



Protein Folding



Weather Forecasting

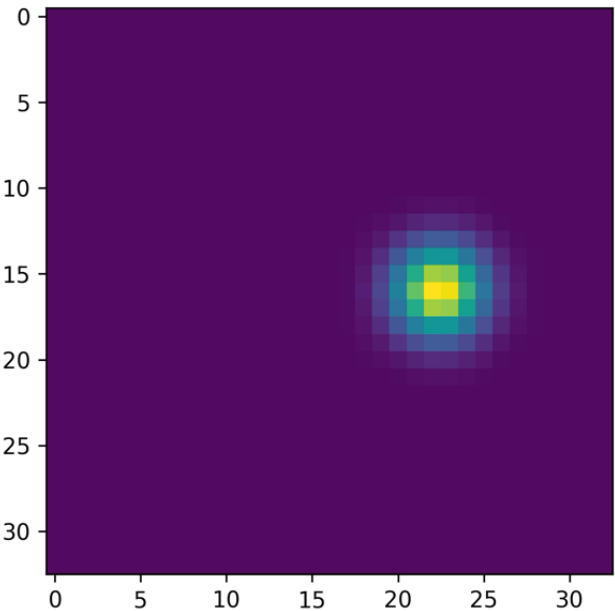


Oil & Gas Exploration Simulation

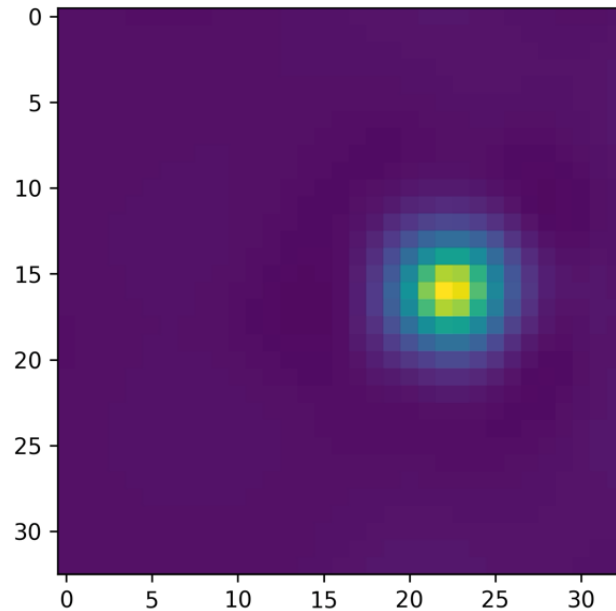
Physics-Informed Neural Networks (PINNs)

Time Step 0 of 151

Solution



PINN



RESULTS ORIGINALLY SHOWN AT SC22, SOLVING A 2D WAVE EQUATION USING A PHYSICS-INFORMED NEURAL NETWORK IMPLEMENTED IN TENSORFLOW 2.

2XIPU FOUND TO BE 11X FASTER THAN 1XA100 GPU, AT SIMILAR MONETARY & ENERGY COST.

WORK DONE IN COLLABORATION WITH STFC HARTREE AND THE UK ATOMIC ENERGY AUTHORITY

Platform	Time to Train / seconds (20k Epochs)	Speedup vs GPU
2 Bow IPUs	41	11x
A100 GPU	530	-

HOT TOPICS IN AI RESEARCH



IPU for GNNs

The next frontier in AI/ML

GRAPHCORE

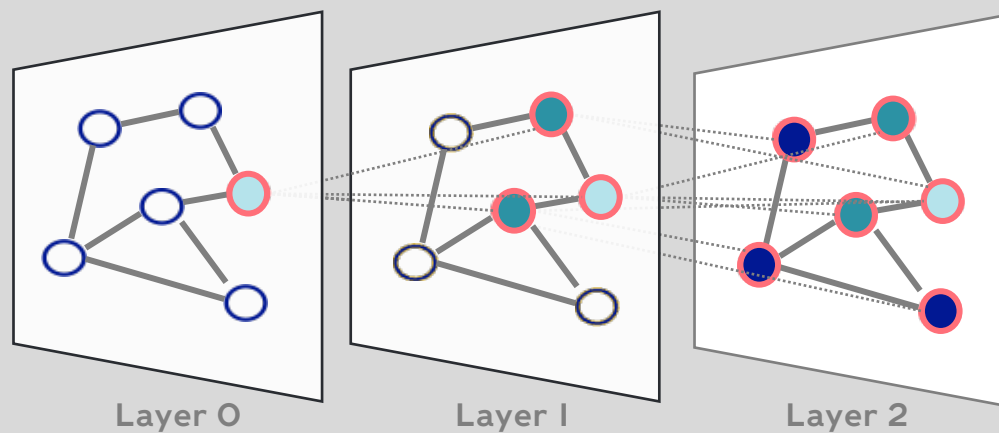




WHAT IS A GNN ?

GRAPH NEURAL NETWORK

GNNs ARE USED TO SOLVE GRAPH PREDICTION TASKS



Visualisation of how a node accumulates information from neighbouring nodes through the layers of the GNN

GNNs broadly follow a recursive neighbourhood aggregation (or message passing) scheme, where each node aggregates feature embeddings of its neighbours to compute its new feature embeddings^[1]

[1] summary derived from 'How Powerful are Graph Neural Networks?' MIT/Stanford - <https://arxiv.org/pdf/1810.00826.pdf>

GNN USE CASES & APPLICATIONS

HEALTHCARE



DISEASE PREDICTION

RNA–disease association
Disease–gene association
COVID-19 spread prediction



DRUG DISCOVERY

Protein structuring
Protein function prediction
Protein / drug interaction
Drug response prediction



MEDICAL IMAGING

Image segmentation
Abnormal detection
Brain connectivity research
Surgical image analysis



PATIENT RISK PREDICTION

Mining EHRs (health records)

INTERNET



SOCIAL NETWORK ANALYSIS

Social influence prediction



RECOMMENDER SYSTEMS

User and item representations



FAKE NEWS DETECTION

Rumor detection + link classification



IDENTITY RESOLUTION

Real-time personalization & advertising

FINANCE



FRAUD DETECTION

Credit card, insurance, loan fraud



TRADE MARKET PREDICTION

Trader connection surveillance



RISK & COMPLIANCE

Risk analytics + compliance reporting



DATA MIGRATION

Data mapping and consolidation



INTEREST RATE RISK

Leveraging credit scores, employment, income, and other socio-economic factors

SCIENTIFIC RESEARCH



PARTICLE PHYSICS

Particle physics simulation



CHEMICAL PHYSICS

QSOR (structure-odor) modeling

TELECOMMS



WIRELESS COMMUNICATION

Power control
Resource allocation
Channel control
Link scheduling

GOVERNMENT



CRIME PREVENTION

Predicting crime associations



FRAUD DETECTION

Anti-money laundering & tax fraud



CONTACT TRACING

Disease contact tracing

TRANSPORT



TRAFFIC FORECASTING

Traffic speed/time prediction

GAMING



FRAUD DETECTION

Collusion detection in gaming



RECOMMENDER SYSTEMS

Online game recommendations

MANUFACTURING



BILL OF MATERIALS

360 degree BOM analysis



TRACEABILITY

Product recall tracing e.g. cars



MASTER DATA MGMT

RDF graph data modelling

SchNet GNN

Modelling Quantum Interactions in Molecules



 Try on Paperspace

Graphcore engineers successfully trained the **SchNet¹** model on IPU's on the **500k water clusters** dataset², to predict the **potential energy per cluster**.

Preliminary results show a time-to-train of **98 minutes** on 2xIPU-M2000, compared with **>60 hours** on 4xV100 GPUs in PNNL's original paper².

¹ JK. T. Schütt¹, H. E. Sauceda, P.-J. Kindermans, A. Tkatchenko, and K.-R. Müller. "SchNet – A deep learning architecture for molecules and materials" *J. Chem. Phys.* **148**, 241722 (2018).

² Jenna A. Bilbrey, Joseph P. Heindel, Malachi Schram, Pradipta Bandyopadhyay, Sotiris S. Xantheas, and Sutanay Choudhury. "A look inside the black box: Using graph-theoretical descriptors to interpret a Continuous-Filter Convolutional Neural Network (CF-CNN) trained on the global and local minimum energy structures of neutral water clusters" *J. Chem. Phys.* **153**, 024302 (2020).



IPU FOR FOUNDATION MODELS

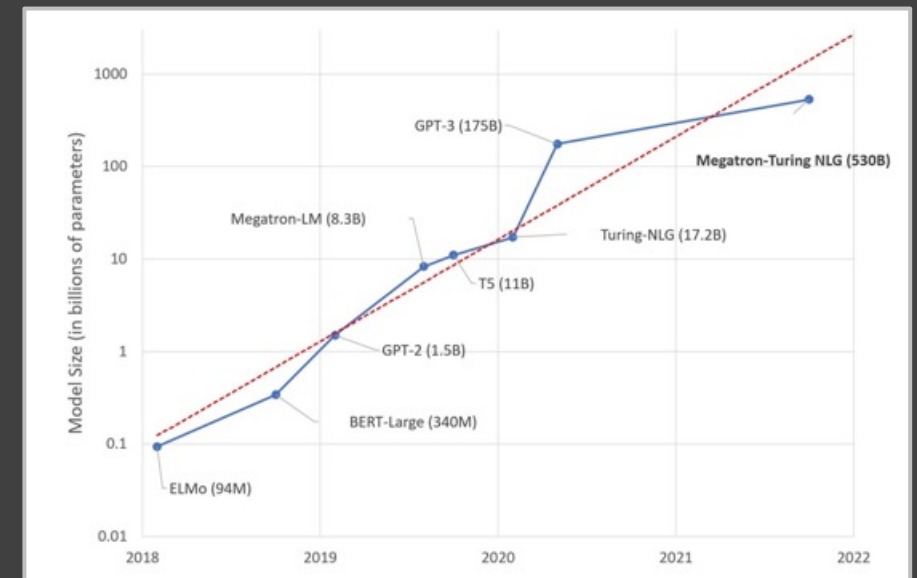


GRAPHCORE



FOUNDATION MODEL TRENDS

- Models are getting much bigger to deliver ever higher demands on improved accuracy & performance
 - This growth is exponential for dense models
- Multimodal models broaden the learning capability by incorporating different modalities (e.g. linguistic, visual, aural)
 - => larger model demands
- Larger dense models mean more compute, more power, more cost
- Counter to this are economic and societal drivers to reduce energy consumption & cost



Exponential trend of SOTA NLP models:

Source: Microsoft/NVIDIA <https://arxiv.org/abs/2201.11990>



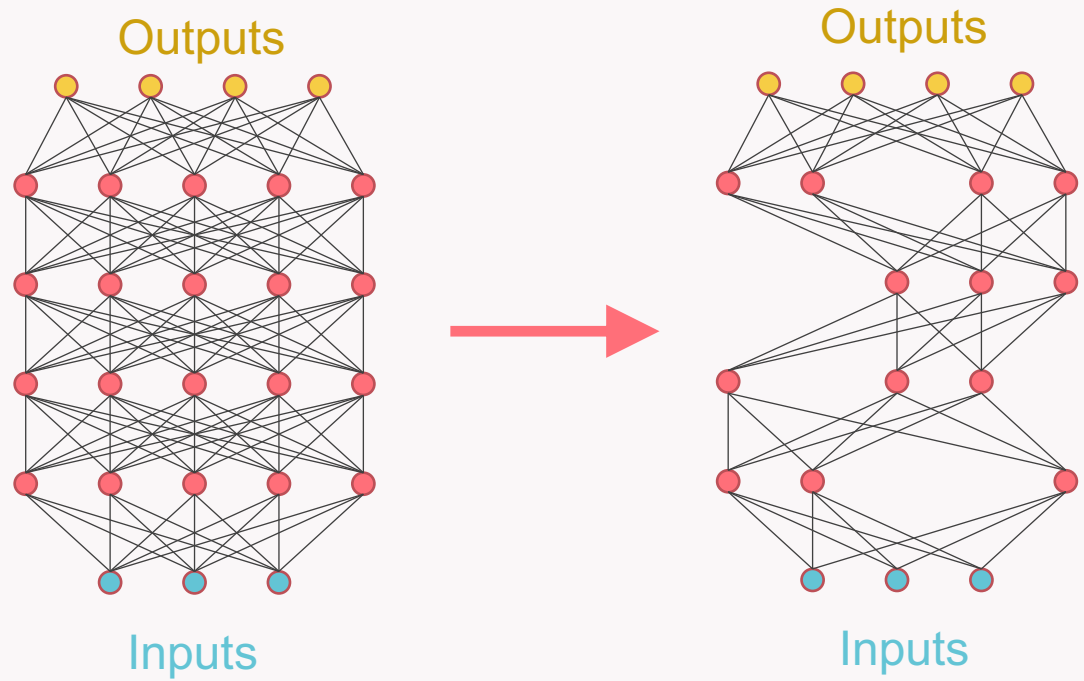
IMPROVING MODEL EFFICIENCY

- **Selectivity / Conditional Models**
 - Models need to become **selective** (or conditional), such as Mixture of Experts (MoE) based models
 - Different parts of models are only used when needed
 - This can help reduce compute growth to linear instead of exponential
- **Sparsification of models**
 - Only incur cost of compute when required
 - Lower memory requirement
 - Fewer multiplications
 - Lower power





WHAT IS SPARSITY?



Only a small proportion of connections are key to model behaviour

So we can prune and re-train to create a “sparse” model

This is beneficial on a processor like the IPU that can do the sparse computation efficiently





THE 'GOOD' COMPUTER

GRAPHCORE



ROADMAP TO ULTRA-INTELLIGENCE AI

Human brain has around 100 billion neurons

With 100Tn+ synapses, equivalent to parameters in an AI model

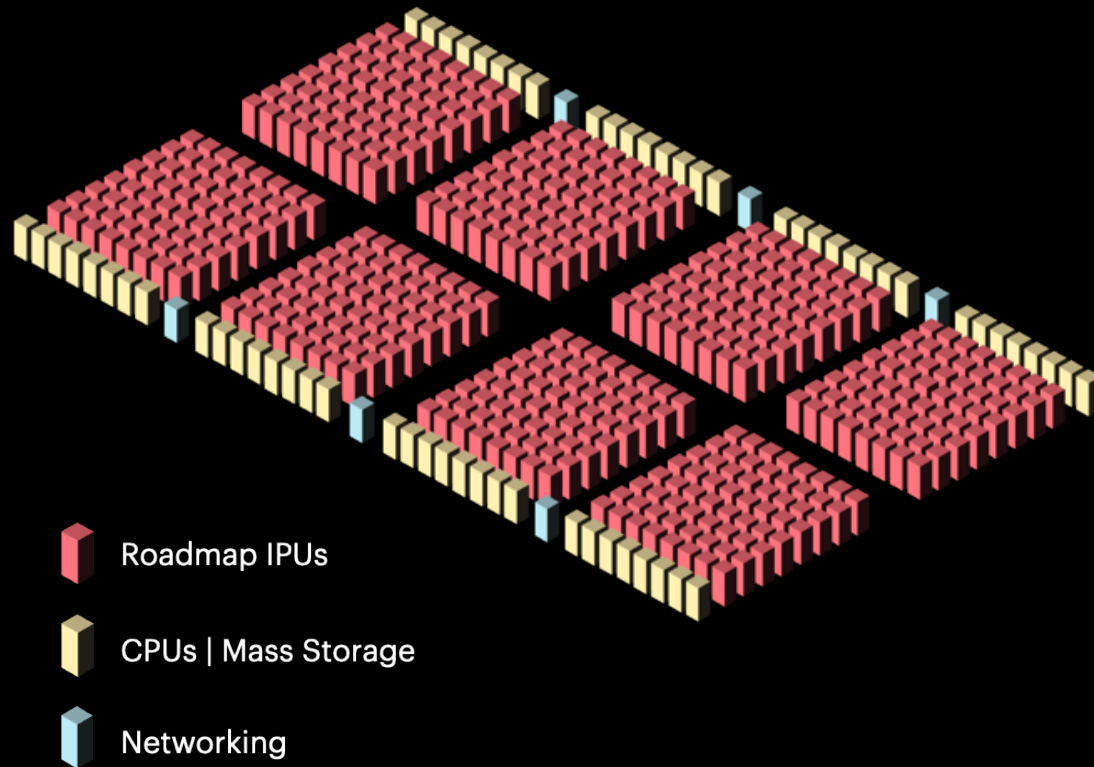
Current largest AI models are around 1Tn parameters

Graphcore is developing an Ultra-Intelligence Machine that will surpass the parametric capacity of the brain

GRAPHCORE



THE 'GOOD' COMPUTER



Over 10 **Exa-Flops** of AI floating point compute from 8,192 roadmap IPUs

3D Wafer-on-Wafer logic stack

Up to **4 PB** of memory with bandwidth of over **10 PB/s**

Enabling AI models to be developed with **500 Tn parameters**

Fully supported by Poplar[®] SDK

GRAPHCORE



APPLICATIONS

Very AI Large-Models:

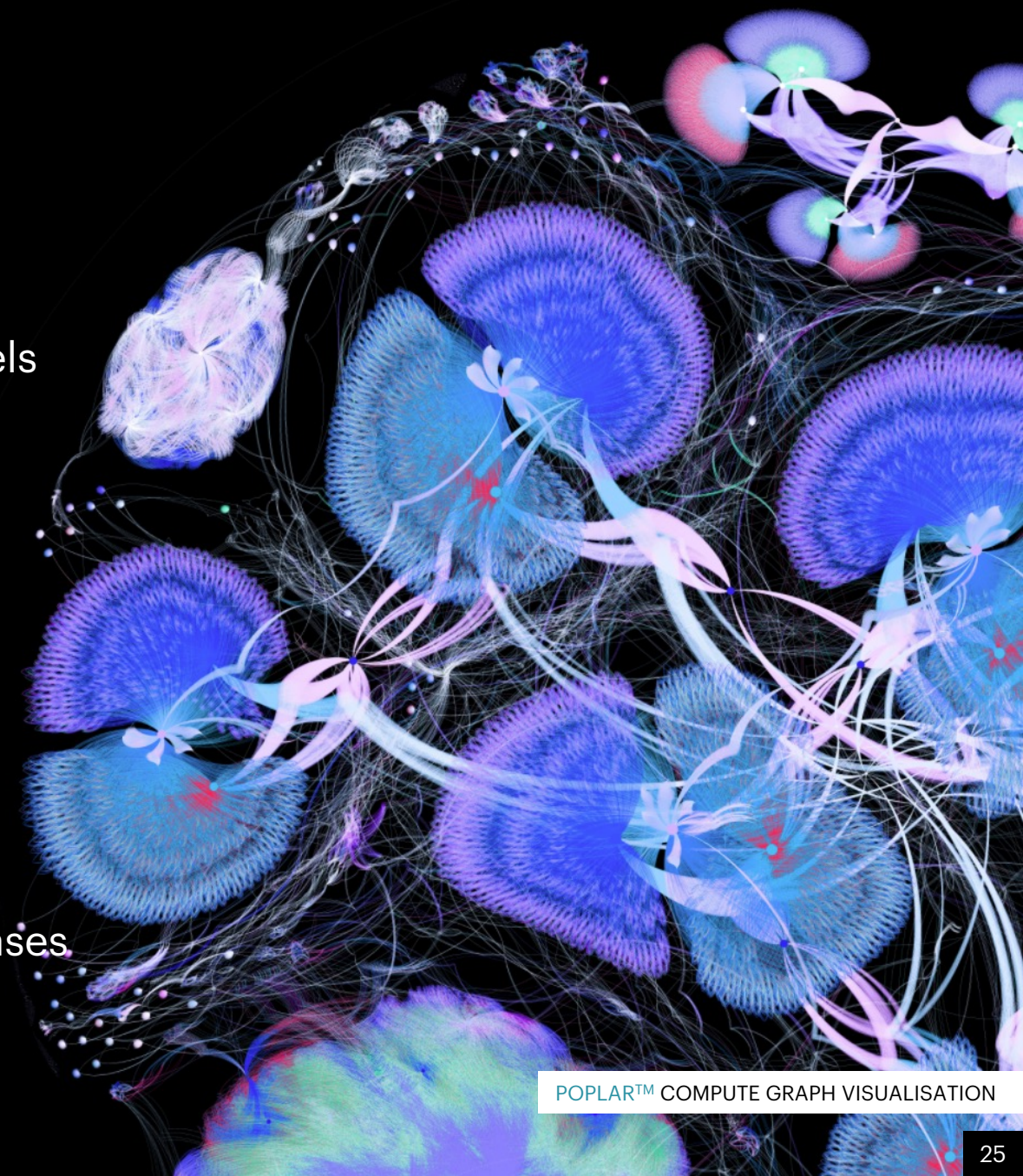
- Multi-trillion parameter model training and inference
- Next-generation, large, conditional and sparse models

AI in Science and Industry

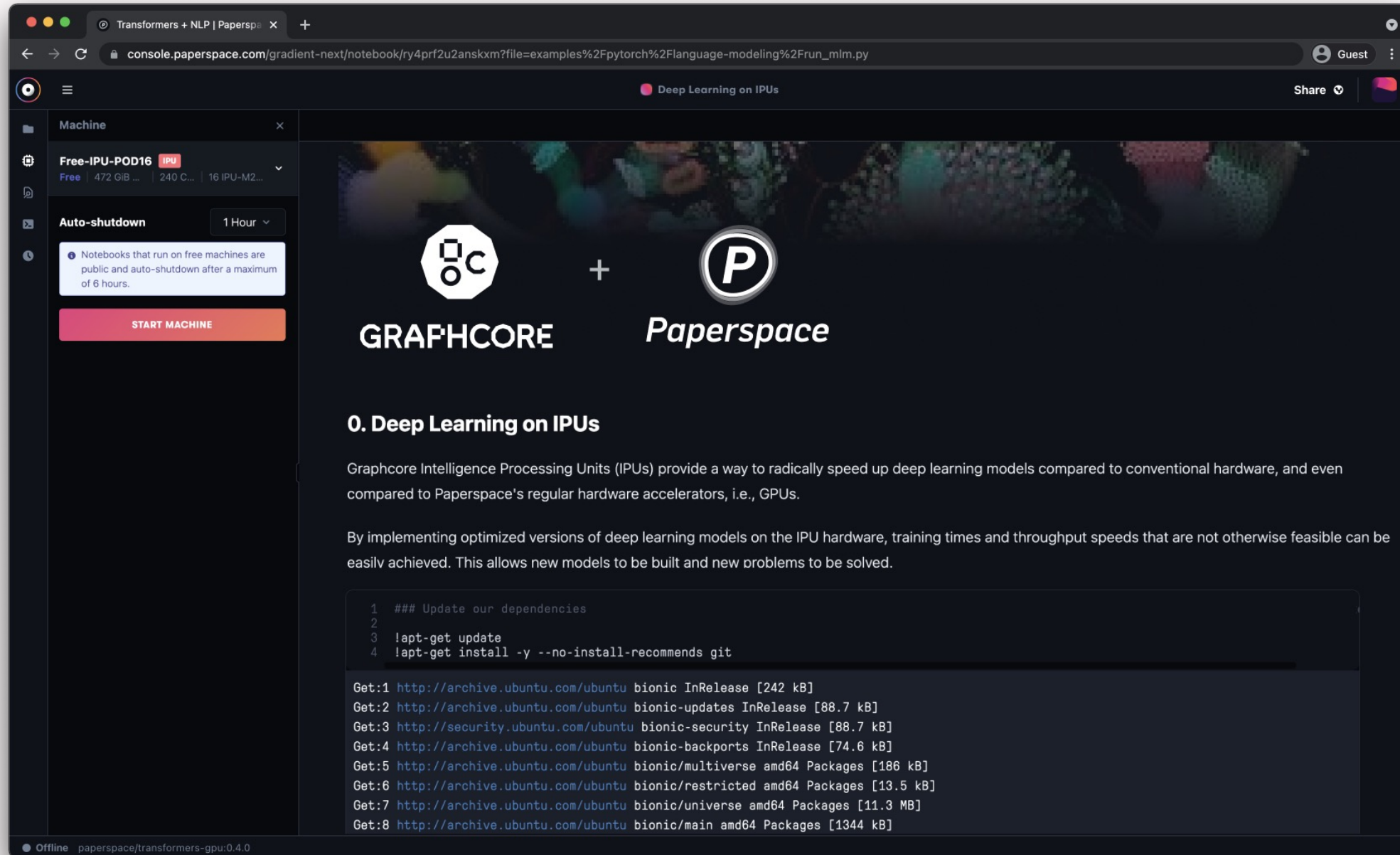
- Healthcare: Genomics | Proteomics | Analysis
- AI-HPC: Simulation | Modeling
- Autonomous systems
- Materials Science | Manufacturing
- Environment: Weather prediction | Smart city

AI in Business

- Language understanding | Process automation | Bots
- Advanced big-data graph analytics and graph databases
- Next generation Recommenders



IPUS IN THE CLOUD



The screenshot shows a web browser window with the URL `console.paperspace.com/gradient-next/notebook/ry4prf2u2anskxm?file=examples%2Fpytorch%2Flanguage-modeling%2Frun_mlm.py`. The page title is "Deep Learning on IPUs". On the left sidebar, the machine configuration is "Free-IPU-POD16 IPU" with 472 GiB of memory and 16 IPU-M2 units. An auto-shutdown timer is set to 1 hour. A "START MACHINE" button is visible. The main content area features the Graphcore and Paperspace logos with the text "0. Deep Learning on IPUs". Below this, there is a paragraph explaining that Graphcore IPU hardware speeds up deep learning models compared to GPUs. A code block contains terminal output for updating dependencies and installing git.

Machine
Free-IPU-POD16 IPU
Free | 472 GiB ... | 240 C... | 16 IPU-M2...

Auto-shutdown 1 Hour

Notebooks that run on free machines are public and auto-shutdown after a maximum of 6 hours.

START MACHINE

GRAPHCORE + **Paperspace**

0. Deep Learning on IPUs

Graphcore Intelligence Processing Units (IPUs) provide a way to radically speed up deep learning models compared to conventional hardware, and even compared to Paperspace's regular hardware accelerators, i.e., GPUs.

By implementing optimized versions of deep learning models on the IPU hardware, training times and throughput speeds that are not otherwise feasible can be easily achieved. This allows new models to be built and new problems to be solved.

```
1 ### Update our dependencies
2
3 !apt-get update
4 !apt-get install -y --no-install-recommends git
```

```
Get:1 http://archive.ubuntu.com/ubuntu bionic InRelease [242 kB]
Get:2 http://archive.ubuntu.com/ubuntu bionic-updates InRelease [88.7 kB]
Get:3 http://security.ubuntu.com/ubuntu bionic-security InRelease [88.7 kB]
Get:4 http://archive.ubuntu.com/ubuntu bionic-backports InRelease [74.6 kB]
Get:5 http://archive.ubuntu.com/ubuntu bionic/multiverse amd64 Packages [186 kB]
Get:6 http://archive.ubuntu.com/ubuntu bionic/restricted amd64 Packages [13.5 kB]
Get:7 http://archive.ubuntu.com/ubuntu bionic/universe amd64 Packages [11.3 MB]
Get:8 http://archive.ubuntu.com/ubuntu bionic/main amd64 Packages [1344 kB]
```

Offline paperspace/transformers-gpu:0.4.0



Free IPU
Access:



IPUS IN THE CLOUD

The screenshot shows the Paperspace Gradient Notebooks console. At the top, the browser address bar displays `console.paperspace.com/google-saml-prod/projects/p5751qgup2t/notebooks`. The interface includes a navigation menu with options like Notebooks, Workflows, Deployments, Models, Data, and Settings. The main content area features the heading "Notebooks" and a sub-heading "Gradient Notebooks is a web-based Jupyter IDE powered by advanced hardware." Below this, it says "Let's create your first notebook." and provides two buttons: "CREATE" and "SEE DOCUMENTATION". At the bottom, a preview of a notebook instance is shown with the title "Predicting Credit Card Fraud" and controls for "STOP INSTANCE", "RESTART KERNEL", and "RUN ALL".



Paperspace

Free IPU
Access:



GRAPHCORE ACADEMIC PROGRAMME

Apply at:
graphcore.ai/academic

Test IPU hardware in
the cloud at no-cost



Support letters for
grants & funding



Access to Poplar® &
PopART® software



Support from
Graphcore Researchers



RESEARCH PRIORITIES:

- Optimisation of Stochastic Learning
- New Efficient Models for Deep Learning & Graph Networks
- Sparse Training
- New Directions for Parallel Training
- Local Parallelism
- Multi-Model Training
- Conditional Sparse Computation

Apply at:
graphcore.ai/academics

GRAPHCORE ACADEMIC PROGRAMMES

Graphcore's Academic Programme provides tools and technology to educators, researchers and students to enable ground-breaking research and innovation in higher education.



MACHINE INTELLIGENCE ACADEMY

The Machine Intelligence Academy is designed to support professors, researchers, and students working in advanced AI and machine learning fields to accelerate their research and develop teaching on IPUs.

[Learn more →](#)



ACCELERATED COMPUTING ACADEMY

The Accelerated Computing Academy is aimed at C++ computer scientists looking to build new applications that transcend today's machine learning, ranging from HPC to simulation modelling, and beyond.

[Learn more →](#)

MACHINE INTELLIGENCE ACADEMY

A new age of IPU-accelerated discovery in computing

WHAT IS IT?

A collaboration initiative designed by Graphcore to enable researchers to develop novel AI techniques and accelerate their research using IPU

WHO IS IT FOR?

Professors, researchers and students working in advanced AI, machine learning and related fields

WHY JOIN?

Members benefit from free IPU cloud credits, support letters, training workshops, engineering support, spotlight promotion and exclusive swag



Access to free IPU hardware in the cloud



Support letters for grant and funding proposals



Bespoke training workshops and educational materials



Support from Graphcore Engineers and SMEs



Project showcase and developer spotlight promotion



ACCELERATED COMPUTING ACADEMY

A new age of IPU-accelerated discovery in computing

WHAT IS IT?

A new computing academy designed by Graphcore to enable new applications that require highly parallel, high-performance compute

WHO IS IT FOR?

C++ computer scientists in academia looking to solve new problems through computationally intensive research that transcend AI and machine learning

WHY JOIN?

Members benefit from free IPU cloud credits, support letters, training workshops, internships, engineering support, spotlight promotion and exclusive swag



Access to free IPU hardware in the cloud



Support letters for grant and funding proposals



Bespoke training workshops and educational materials



Support from Graphcore Engineers and SMEs



Project showcase and developer spotlight promotion



THANK YOU

Dr Alex Titterton
alexandert@graphcore.ai

Academic
programme:



Free 6-hour
IPU Access:





WHY IPUS FOR GNNS

GNN REQUIREMENTS

Low arithmetic intensity due to large memory bandwidth requirements

GNNs often utilise multiple small graphs (or sub-graphs/clusters) & present unique gather/scatter requirements. These require memory intensive operations in parallel

Graphs data structure is highly sparse, hardware capable of handling sparsity efficiently will have an advantage

Developers want to use high-level standard ML frameworks optimised for GNN

Dynamic graphs changing over time require small batch sizes

IPU ADVANTAGE

Ultra-fast, large In-Processor Memory removes memory bandwidth constraints

Truly parallel implementation enabled by IPU's unique MIMD architecture

Fast gather/scatter operation combined with distributed nature of IPU make sparsity its natural domain

Standard ML Framework support including GNN focussed PyTorch Geometric

Optimised small batch size performance