

The past present and future of AI in geosciences

What have we learned and where are we going ?

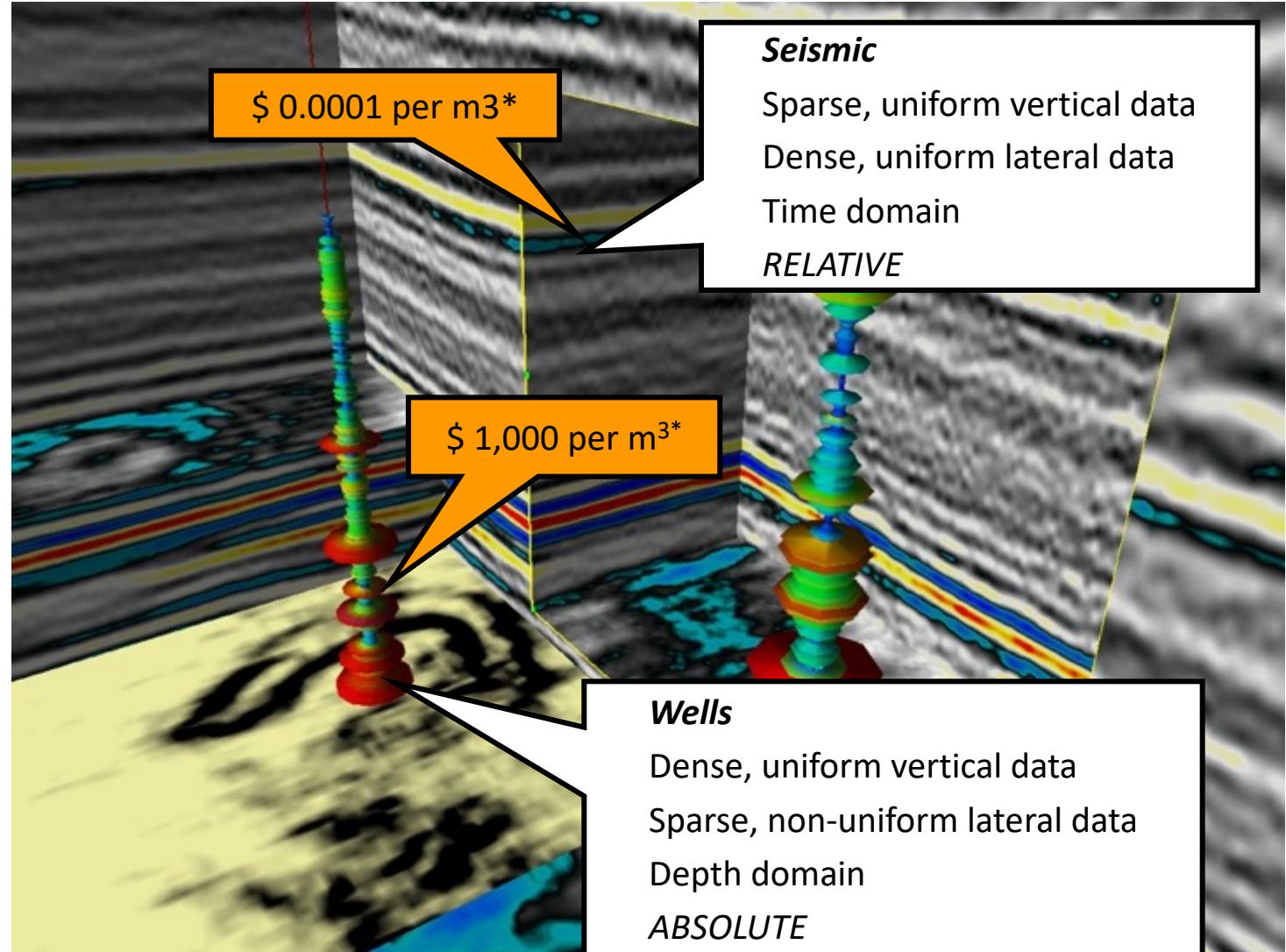
Dr. Lucy MacGregor



Introduction

What data are we dealing with ?

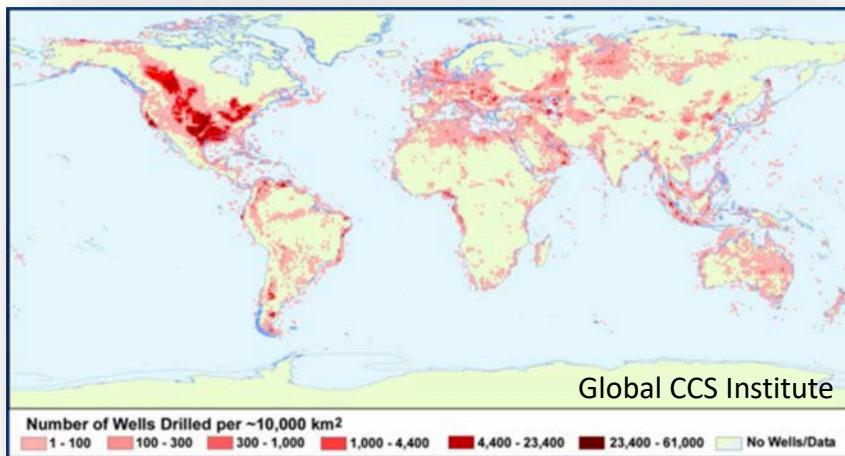
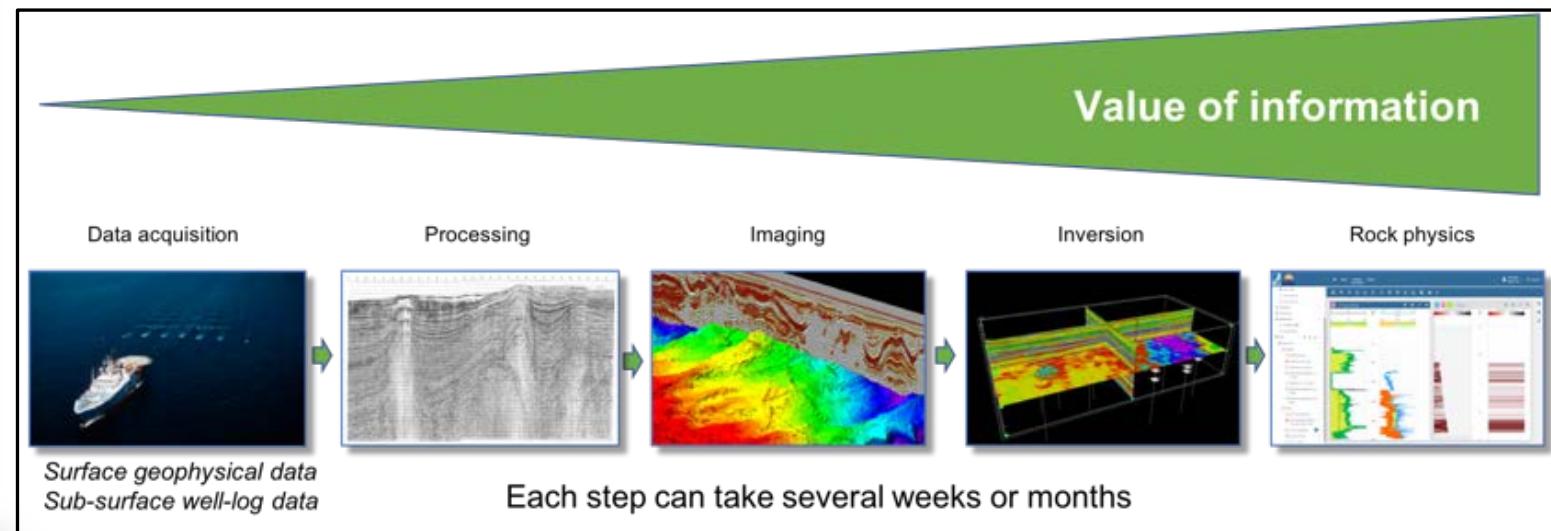
- Seismic data
- Well log data
- Electromagnetic data
- Gravity data
- Magnetics data
- Geological information
- Cores
- Sediment samples
-



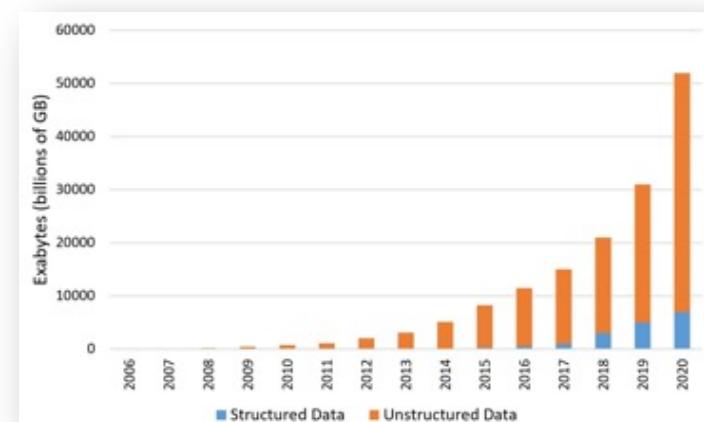
The problem with data:



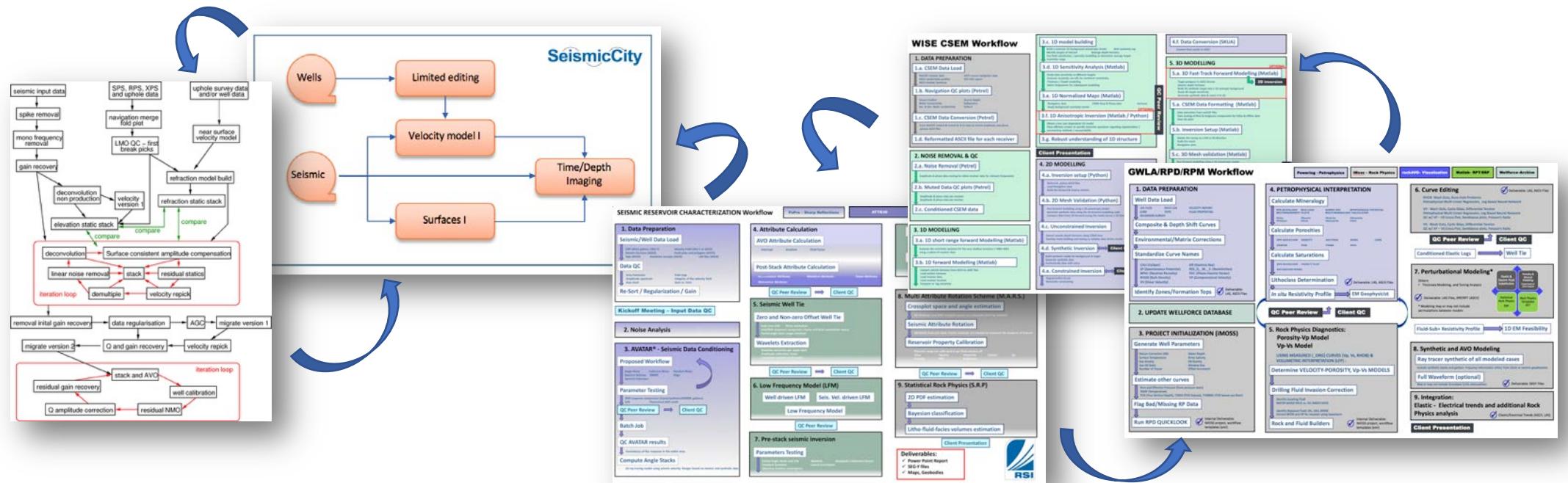
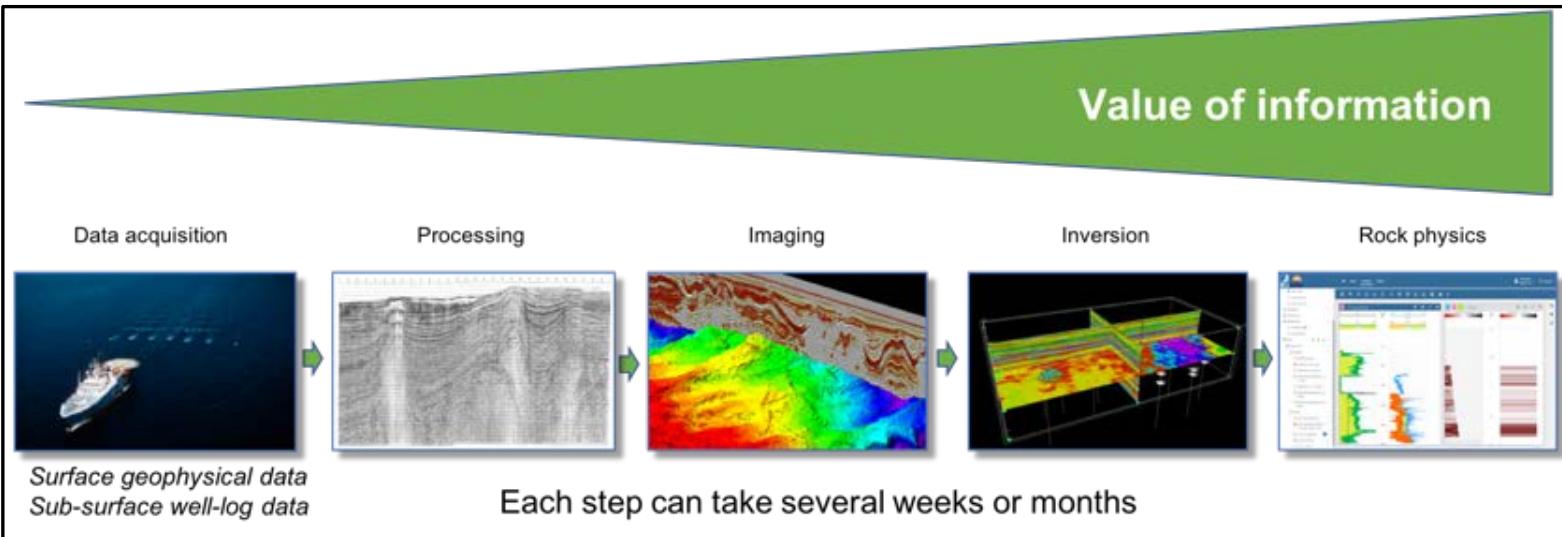
....but it's a bit of a mess...



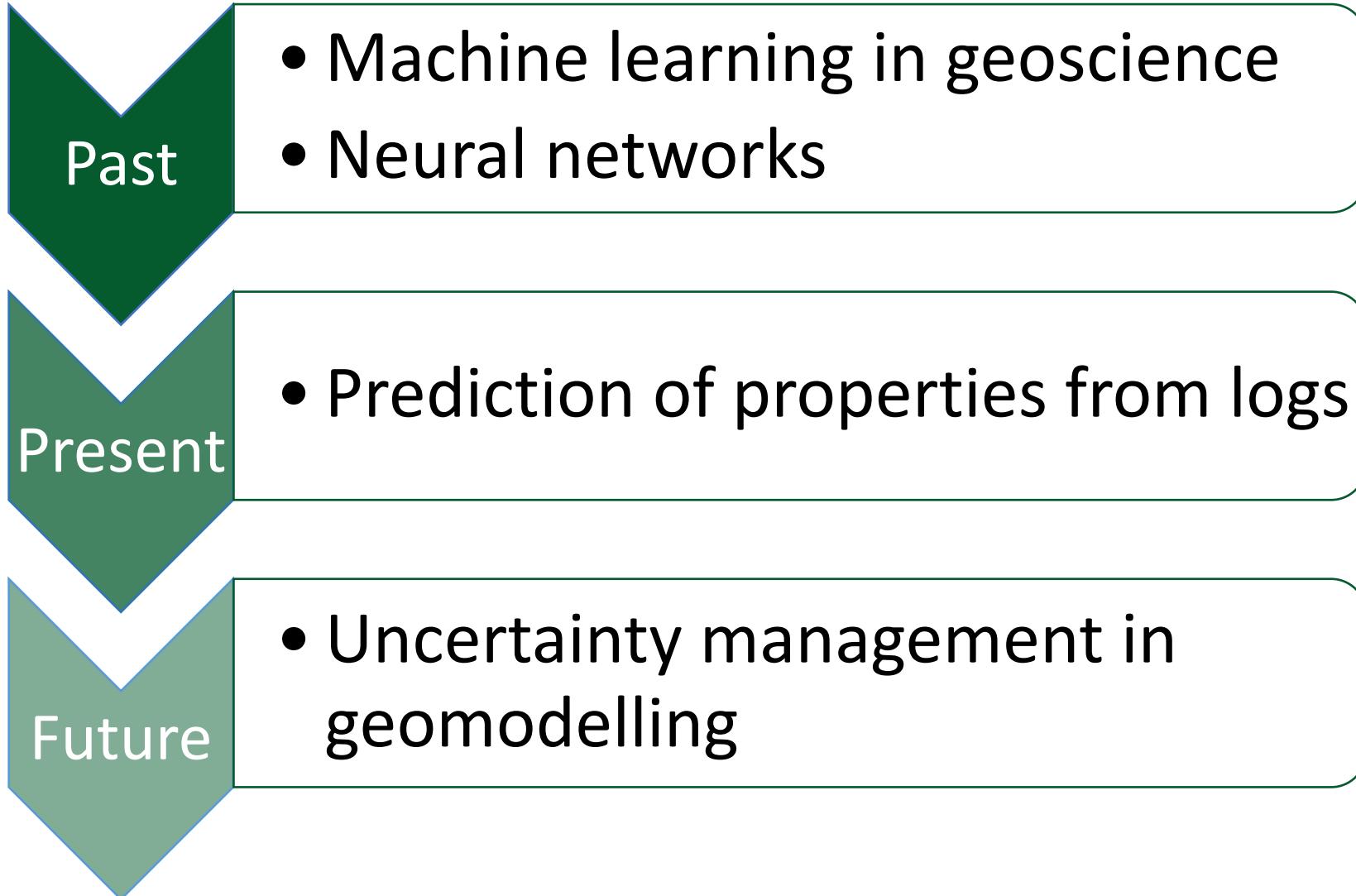
We have a lot of data !



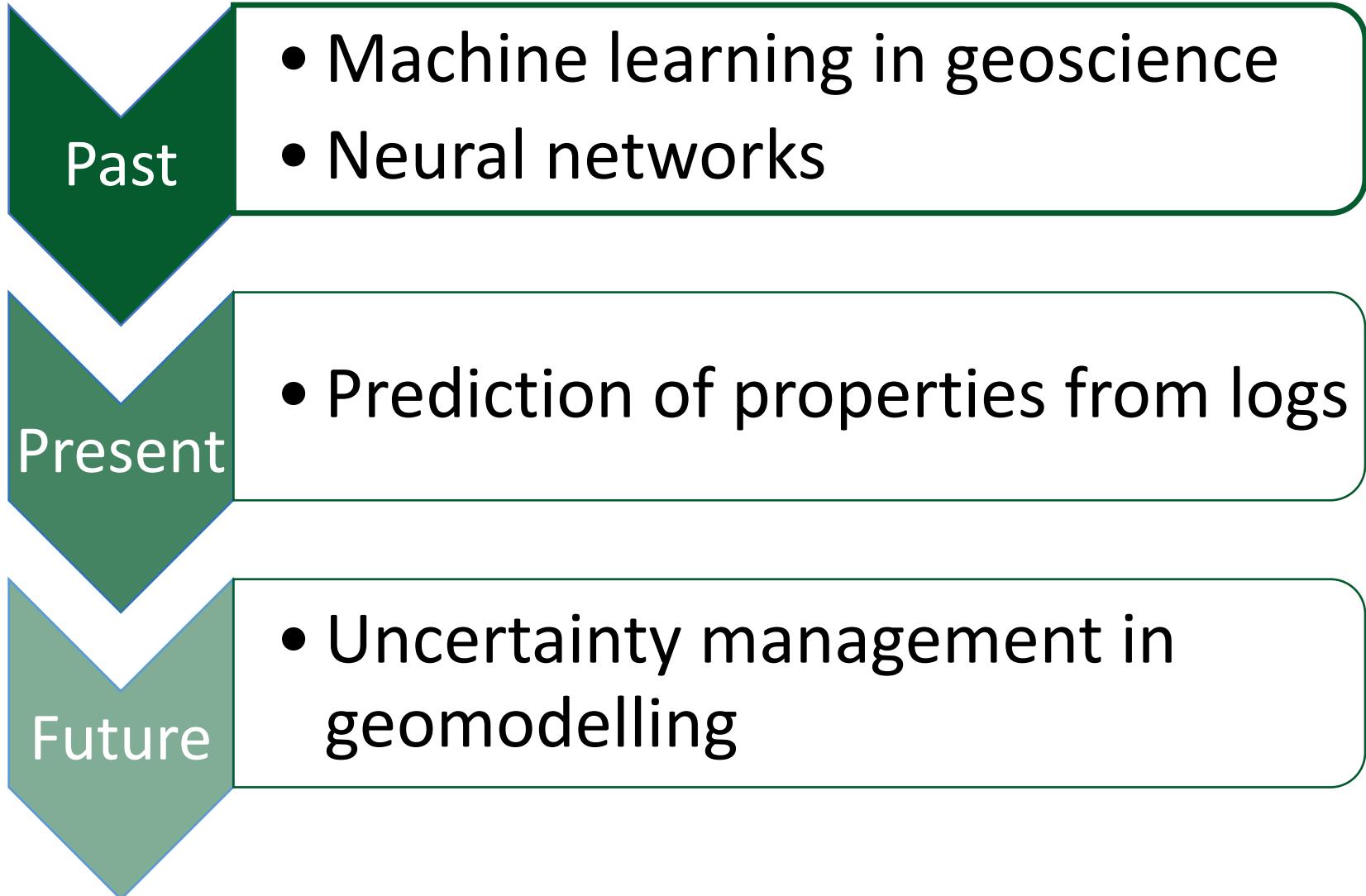
The problem with data:



Overview

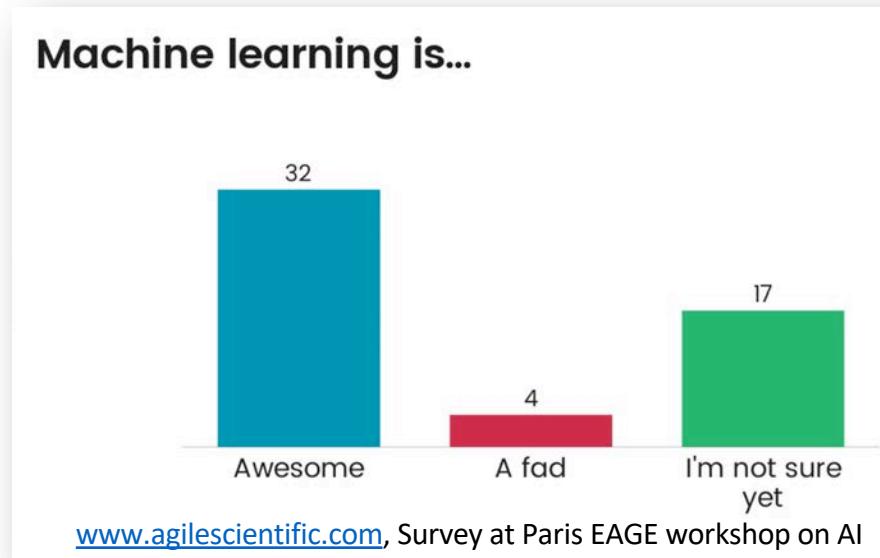


Overview



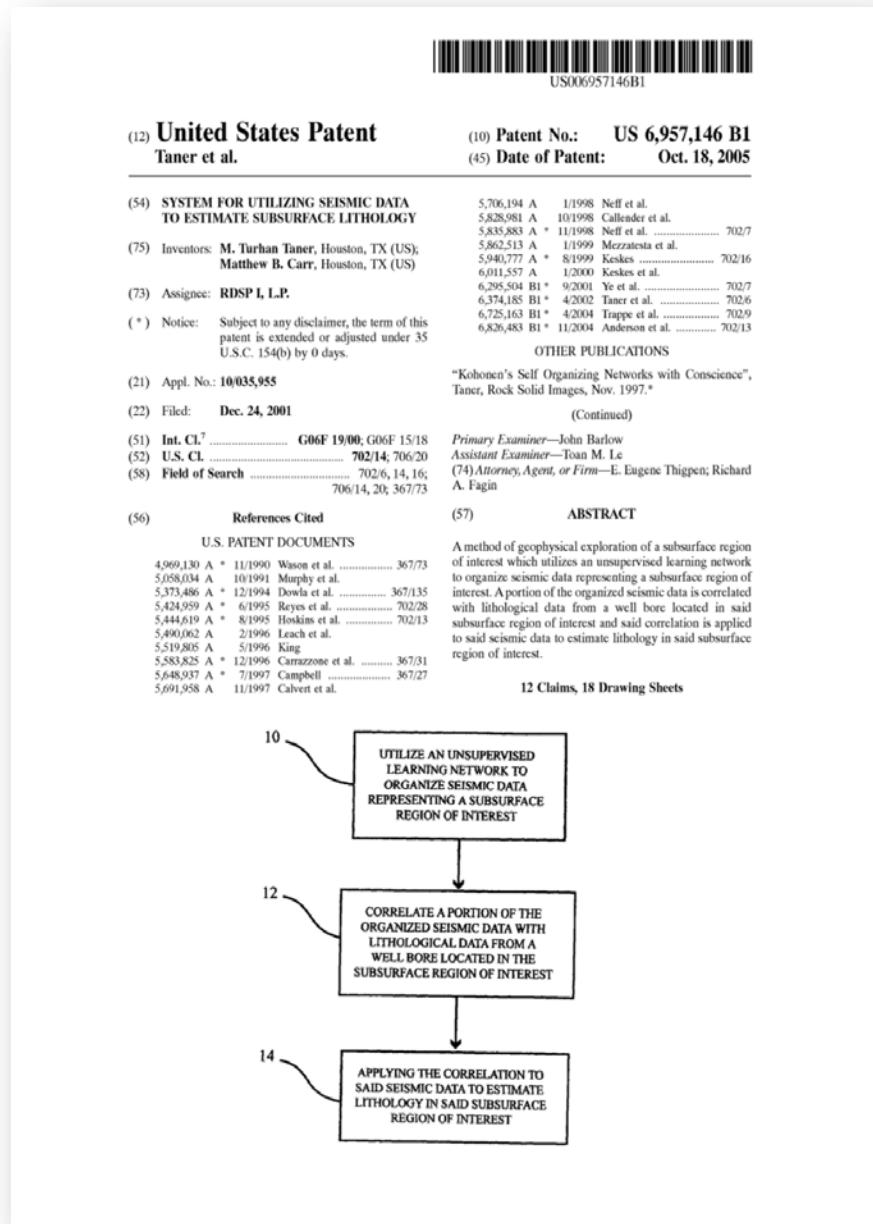
Machine learning in geoscience

- Most of the steps required in geoscience interpretation involve pattern recognition or classification.
- Many machine learning algorithms boil down to a regularized inversion for a function linking predictors and outputs.
- As geoscientists we've inverting things for years !
- However as an oil industry we've been overtaken by e.g. google, amazon....

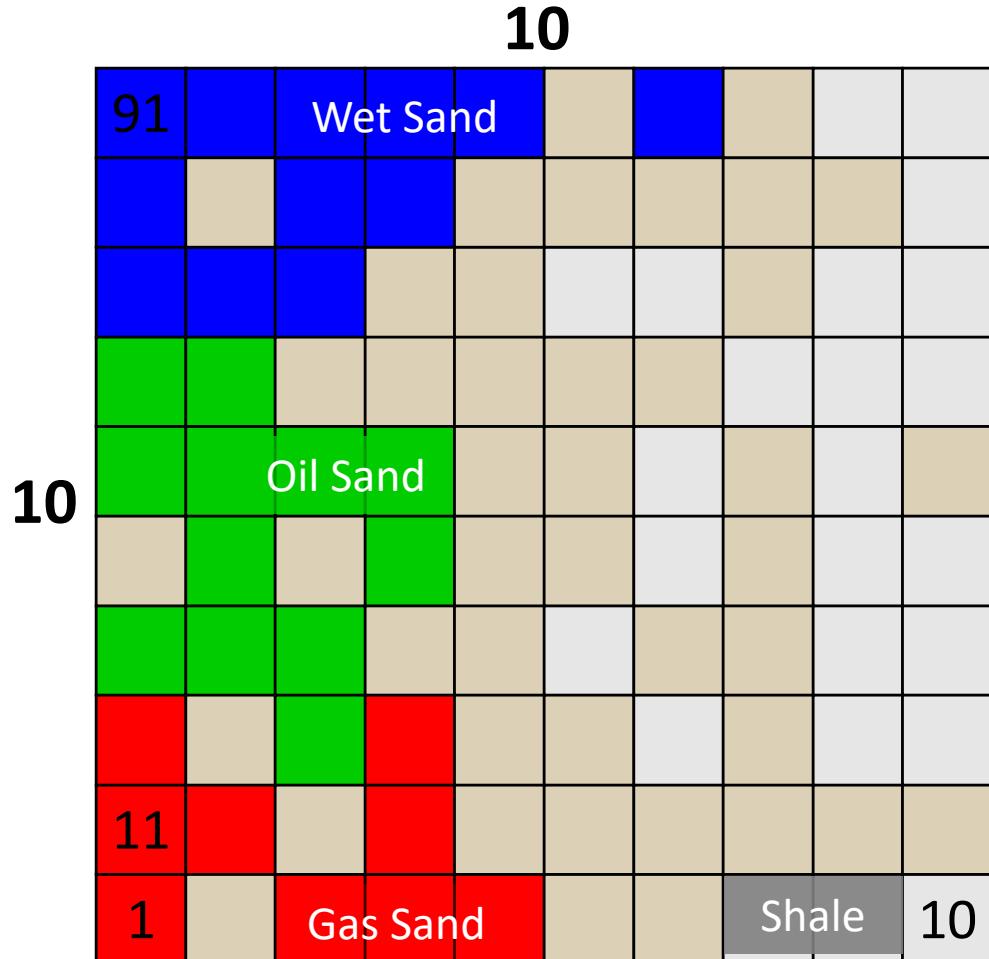


An example: LITHANN®

- Neural network classification of seismic attributes.
- Developed by Tury Taner in the late 1990s as part of RSI's LFP consortium and patented.
- Uses a Kohonen self organizing map
- Still in use today !



How Do KSOM's Work?



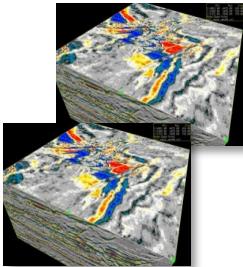
Kohonen self-organizing map algorithm posts ANN classifications of multiple seismic attributes within the confines of a 2D topology

Samples (seismic facies) exhibiting a similar multi-attribute signature are mapped into contiguous regions of Kohonen space.

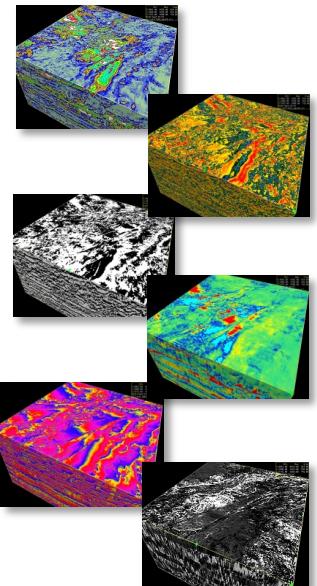
10 X 10
(100 Classes)
2D Network Topology

LITHANN® workflow

Conditioning → Calculation → Classification → Calibration

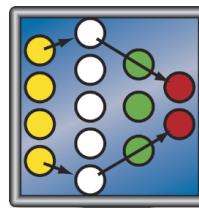


Multiple pre-stack
&/or poststack
seismic volume(s)

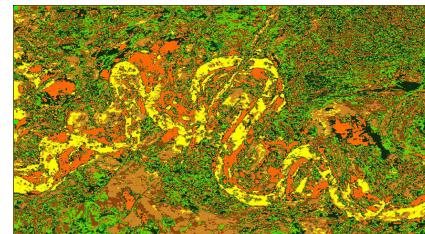


instantaneous, elastic,
geometric, waveshape,
attribute volumes

Classification

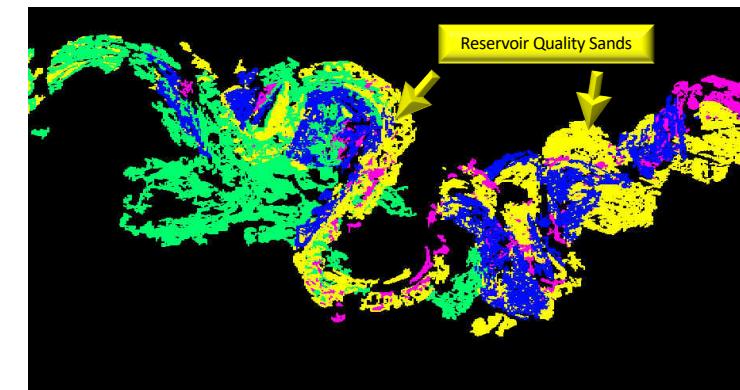


1 classified volume:
2-D topology, 64, 100, 144, or 225
classes
unsupervised or
supervised



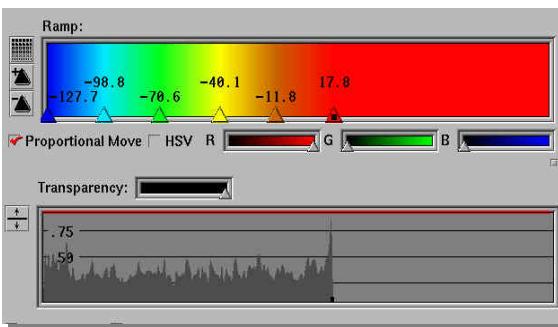
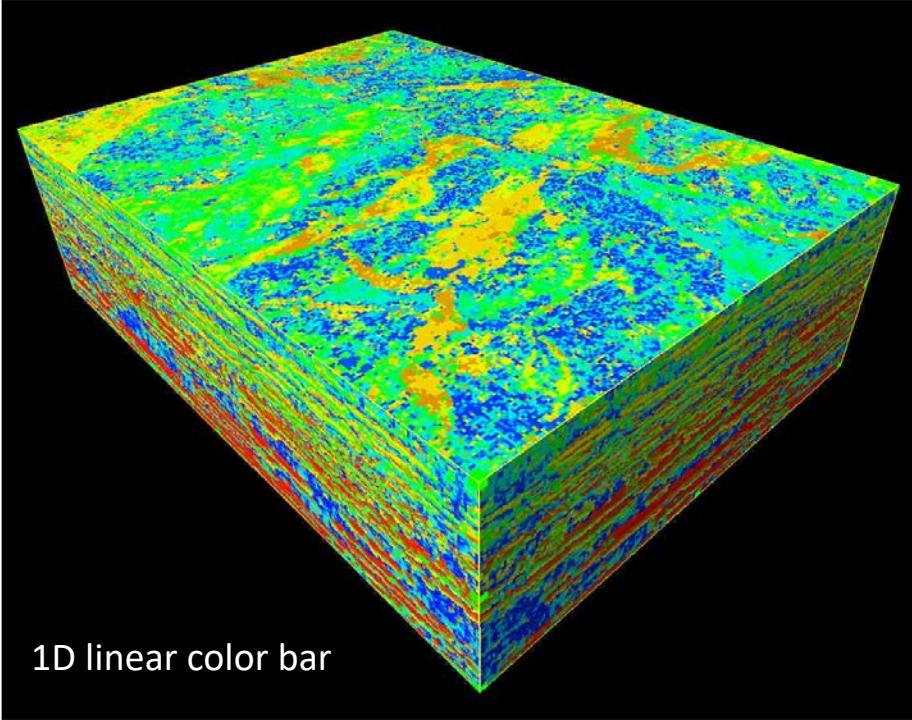
Neural net

Calibration

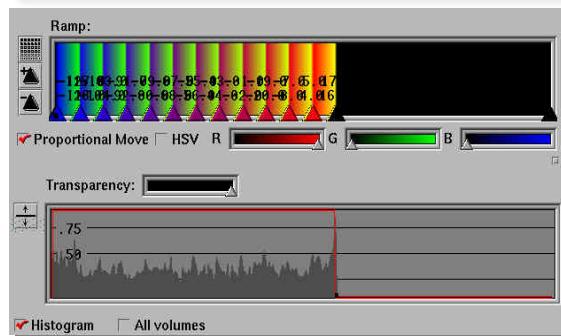
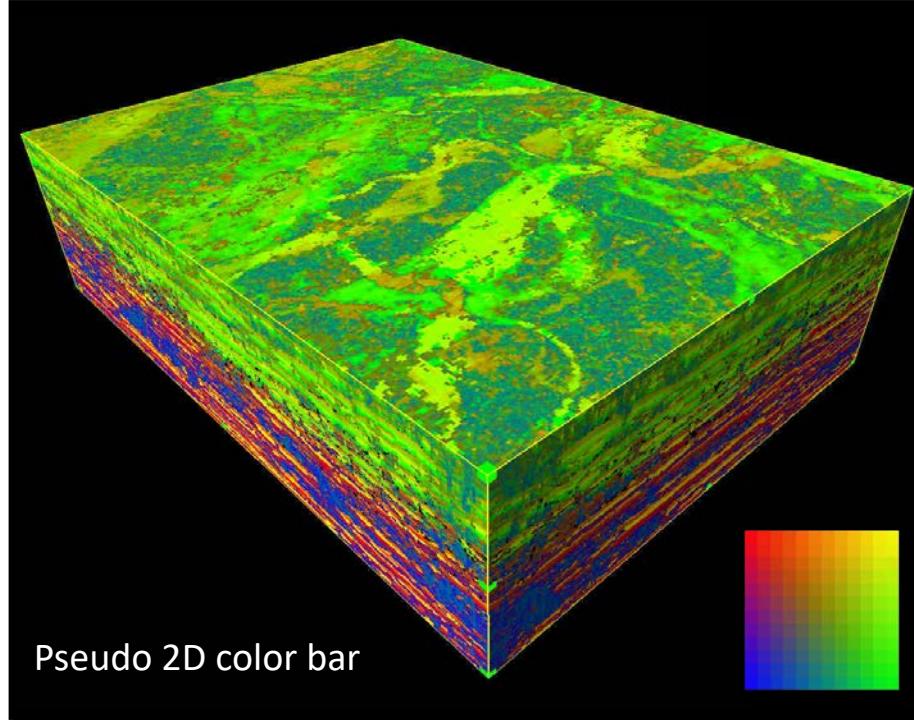


1 seismic facies volume:
3 calibrated classes

Seismic Facies Visualization



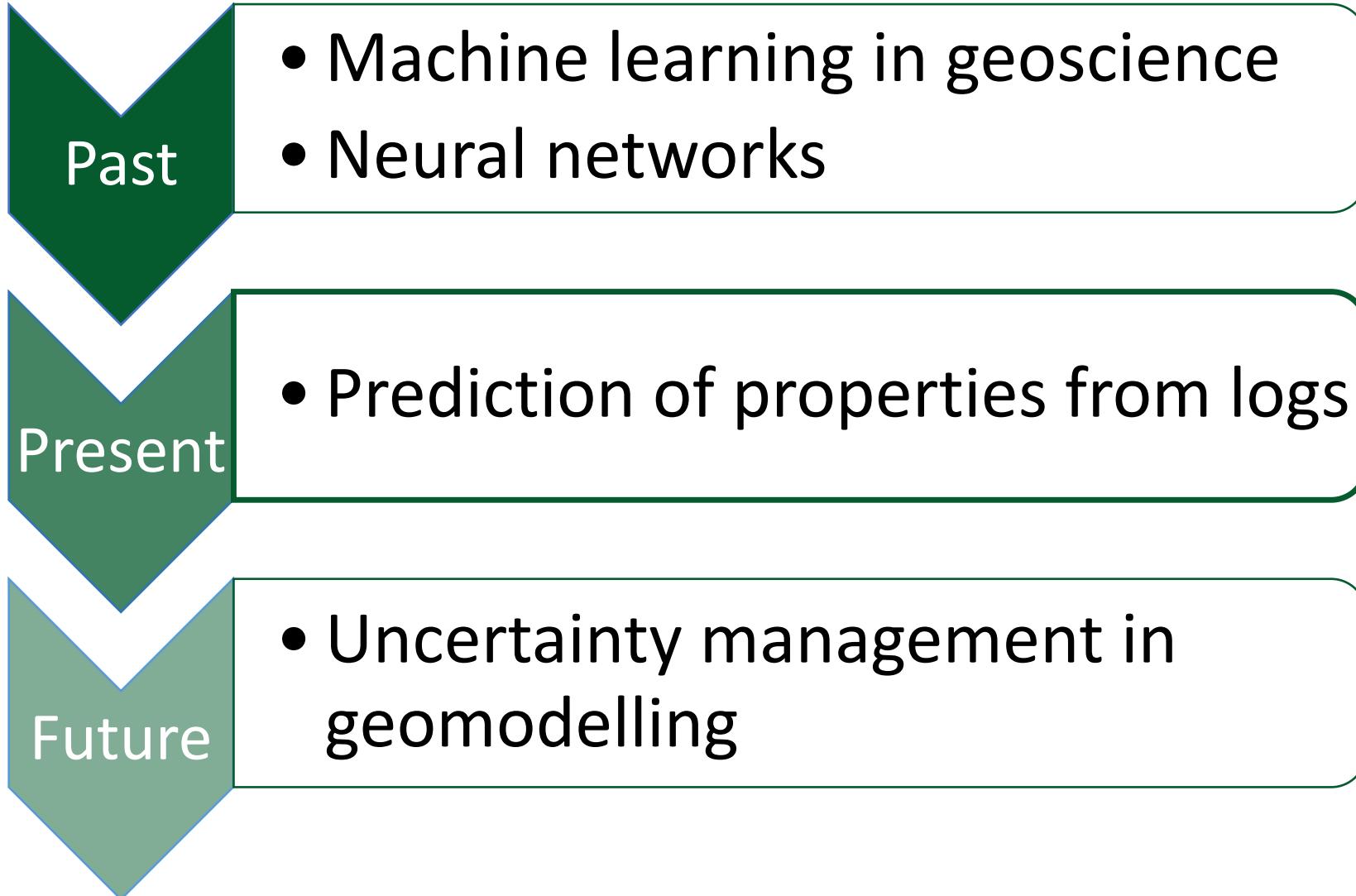
Apparent decrease
in spatial
resolution



Apparent
increase
in spatial
resolution



Overview



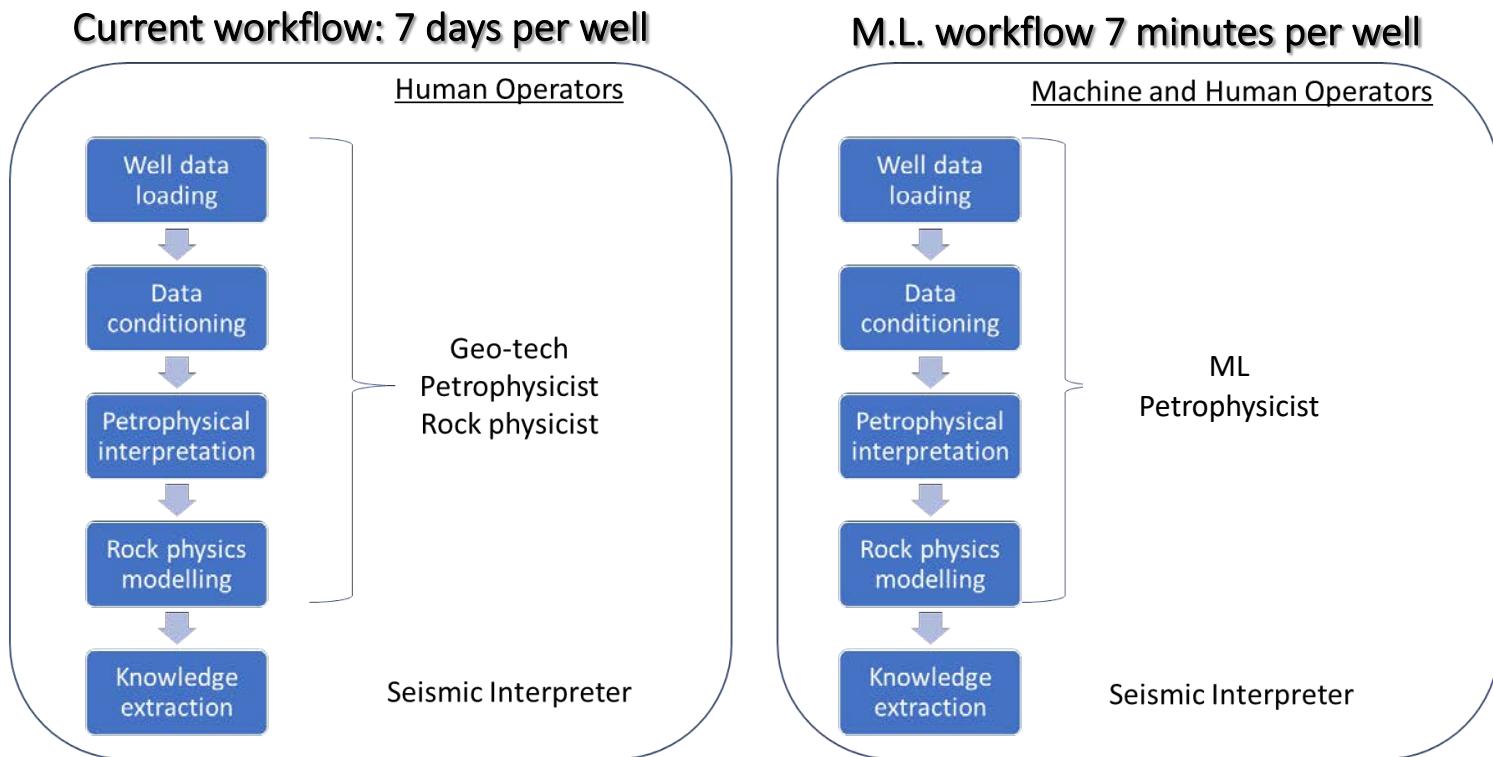
There has been an explosion of ML libraries:

- Most are open source: access to numerous advanced algorithms and a community of users
- (Relatively) easy to use
-but the key is training data



Project SWOOP: Streamlined WOrkflows for Optimised Petrophysics

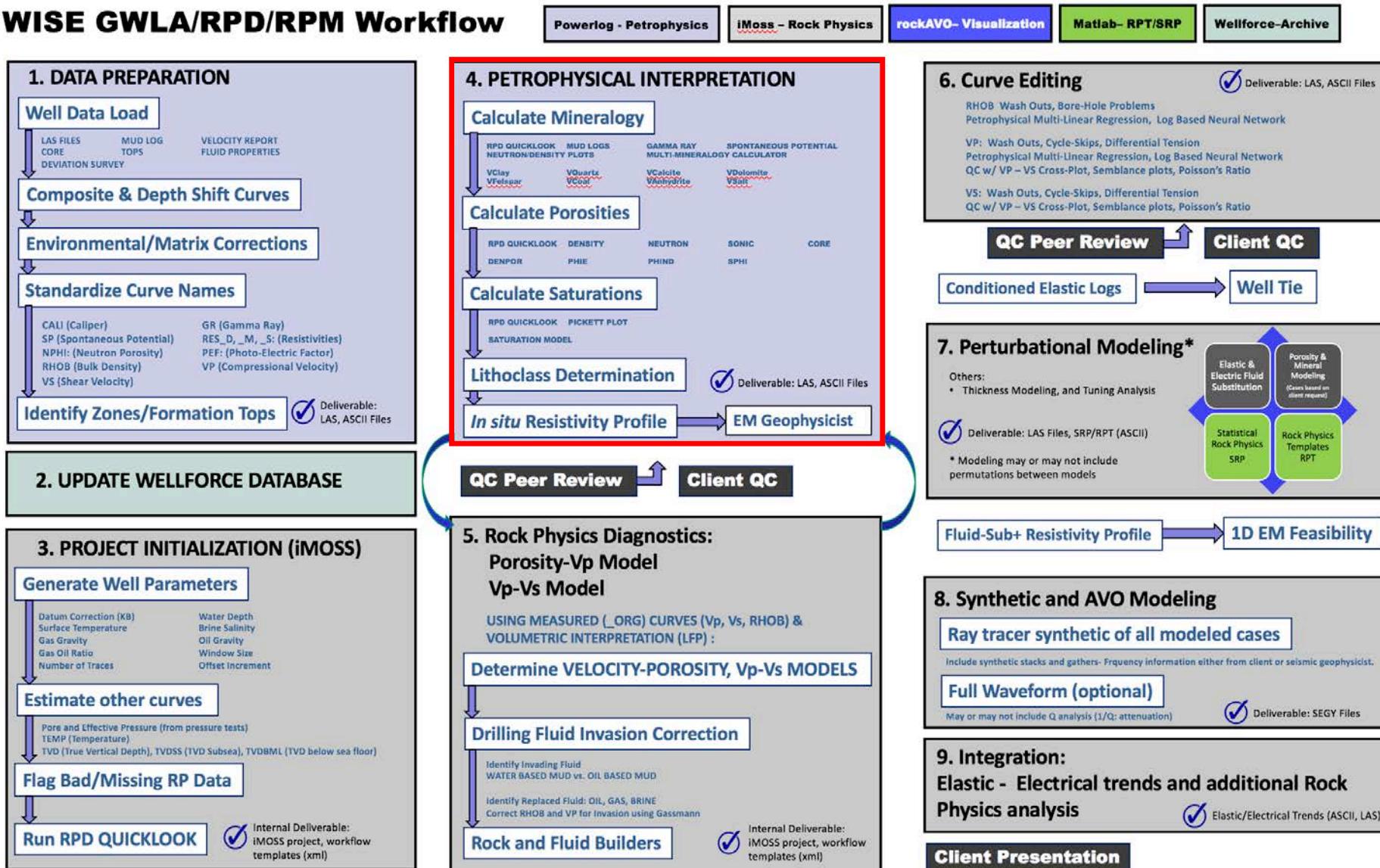
The project was conducted in partnership with EPCC – Edinburgh University's AI team in the high performance computing centre.



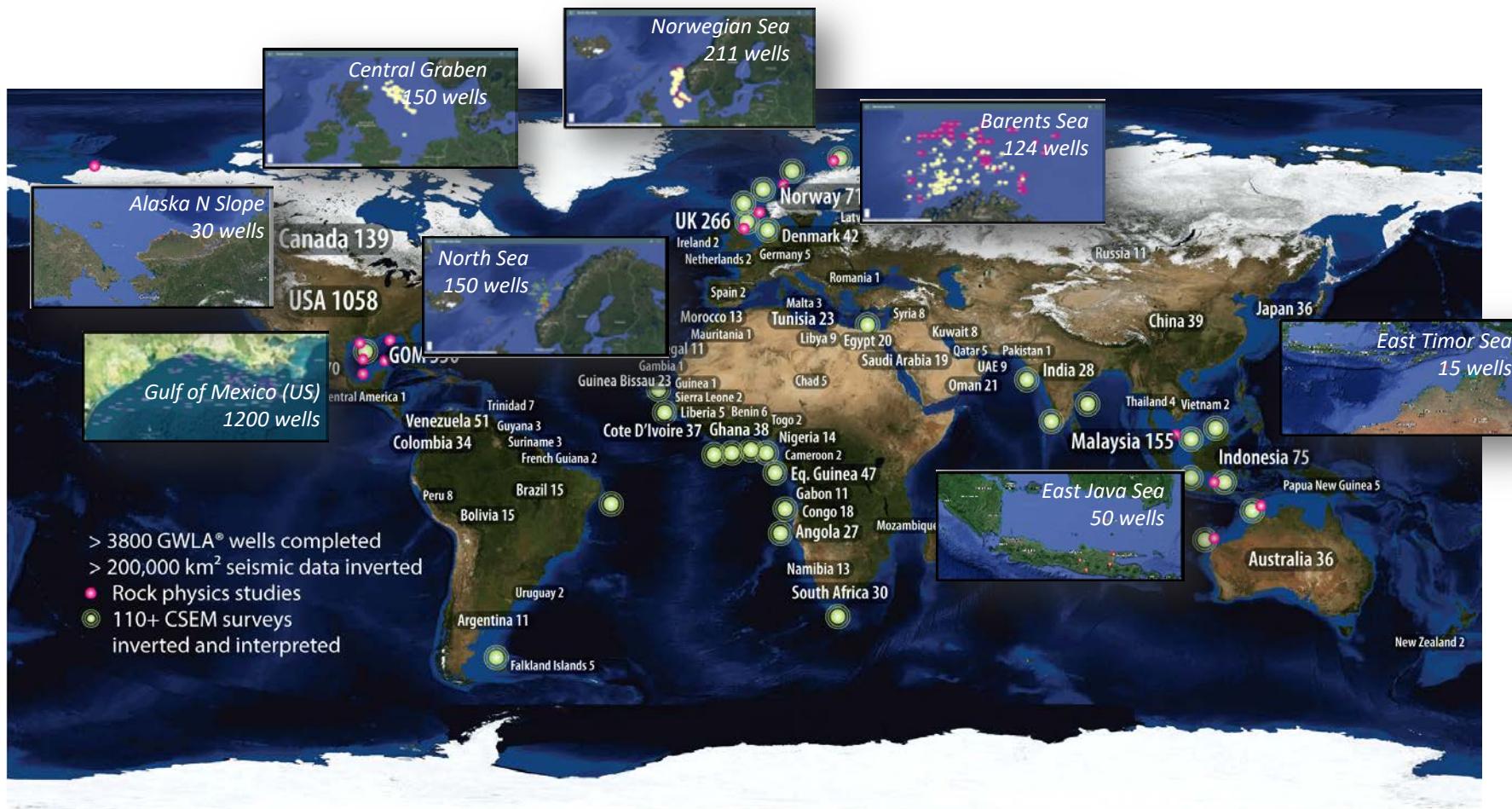
- Streamlining petrophysical workflows with machine learning, MacGregor, L., Brown, N., Roubickova, A., Lampaki, I., Berrizbeitia, J., Ellis, M. & Nichols, K., 2019, EAGE Conference on Reservoir Geoscience, December 2018
- Machine learning on Crays to optimise petrophysical workflows in oil and gas exploration, Brown, N., MacGregor, L., Roubickova, A., Lampaki, I., Ellis, M., Vera de Newton, P., 2019, Journal of Distributed and Parallel Computing, subm.

Petrophysics/rock physics workflow

WISE GWLA/RPD/RPM Workflow



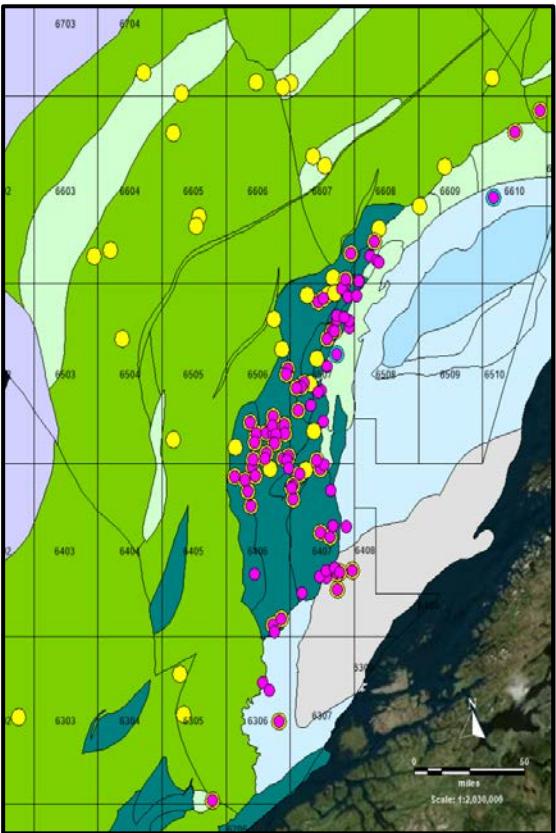
Training data: must be CONSISTENT and COMPLETE



2000 wells in the existing multi-client database

Starting point – training data and ML algorithm

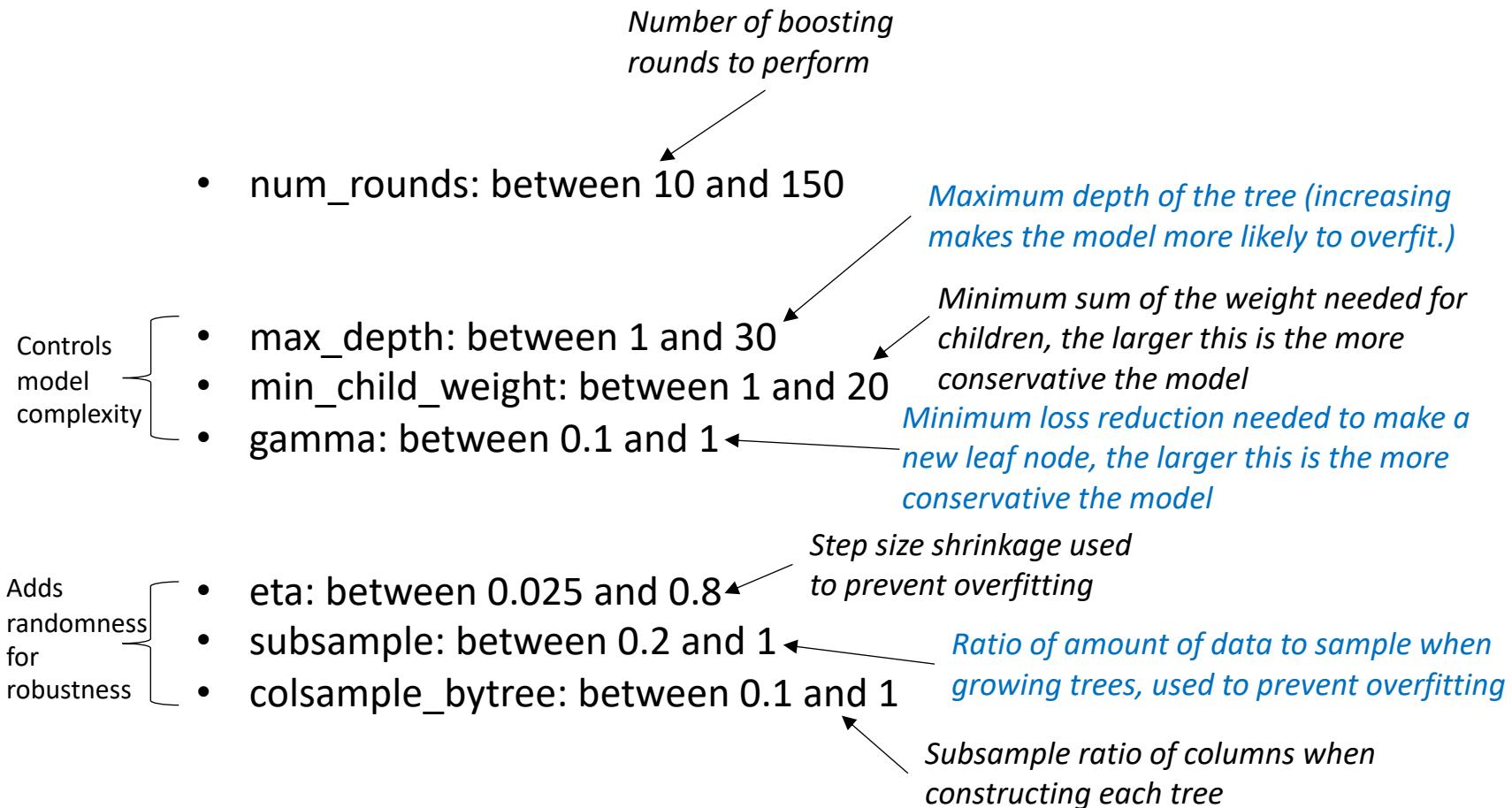
Training data:



64 wells from mid-Norway

- Which algorithm ? We have tried:
 - Dirichlet regression (didn't work very well)
 - Deep Neural Network (current second favourite)
 - Multi-layer perceptron
 - Boosted trees (current favourite)
- How to choose ?
 - Currently looking into a parameter optimisation approach.

Optimising hyperparameters



Clay: num_rounds=142, max_depth=3, min_child_weight=3.5, gamma=0.6, eta=0.05, subsample=0.7, colsample_bytree=0.9

Quartz: num_rounds=120, max_depth=5, min_child_weight=17.5, gamma=0.2, eta=0.15, subsample=1.0, colsample_bytree=0.1

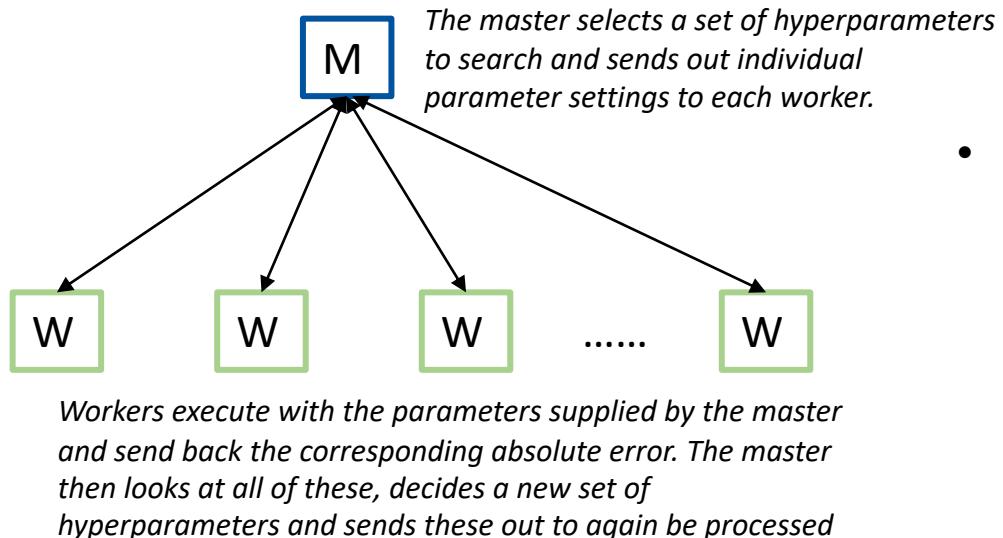
Calcite: num_rounds=135, max_depth=3, min_child_weight=14.5, gamma=1.0, eta=0.075, subsample=0.4, colsample_bytree=0.75

Use hyperopt to search through the parameter space

- Select five different splits between training and test data and for each parameter setting train and test the model with each of these
- Looking to optimise the mean absolute error between prediction and truth here and hyperopt will direct its searching based on these values
- Whilst each of these hyperparameters has a specific meaning, they are interlinked and the best settings are not always obvious (so best to find empirically.)

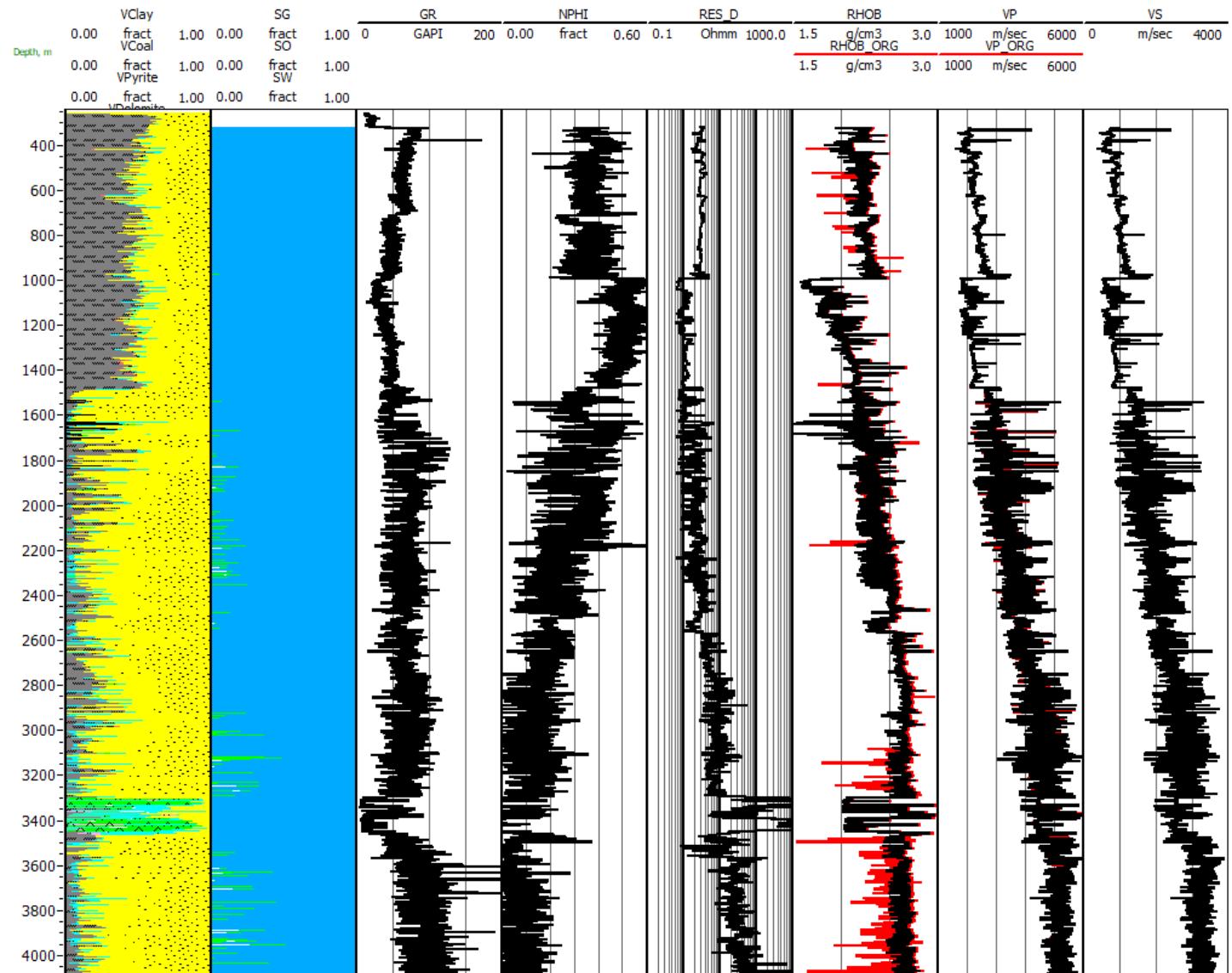
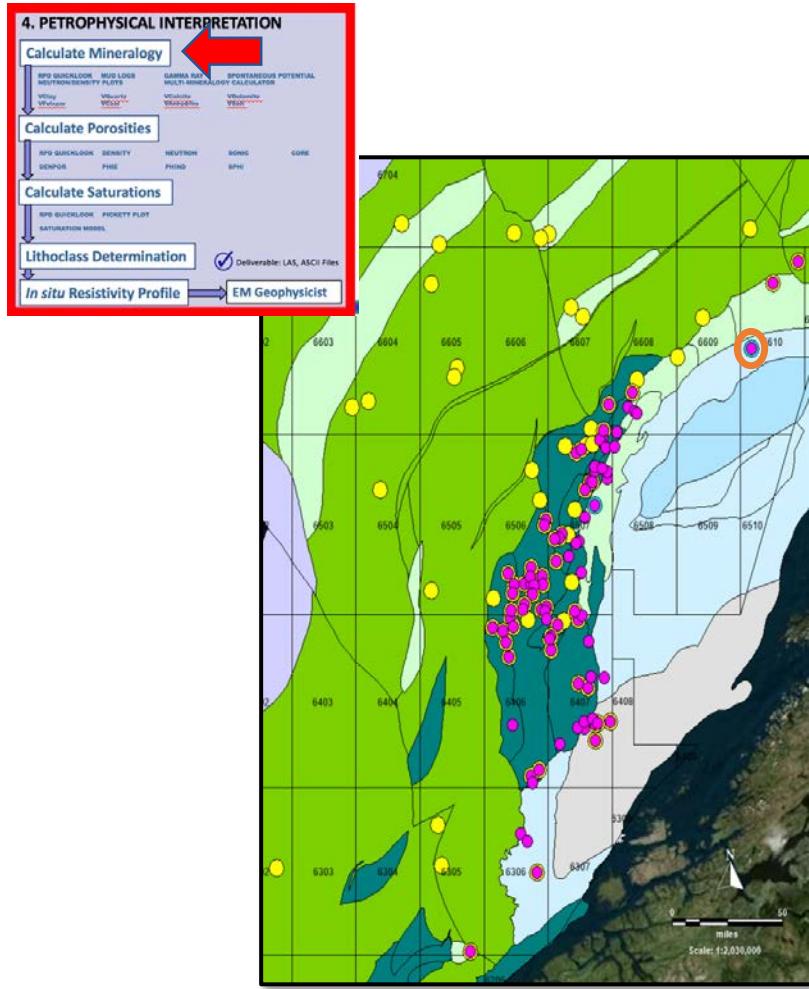
Speeding up hyperparameter searching

- Searching for hyperparameters can take a long time – with the largest number of input curves around 20 minutes per parameter
 - Which means only around 72 parameters can be searched in a day
 - Annoyingly the XGBoost library is not parallelised over multiple CPU cores and the GPU support is quite immature (it tends to run out of memory on our GPU machine.)
- Using Cirrus, an HPC machine and even in a single node there are 36 physical cores. But hyperopt is not able to take advantage of these cores.
- Therefore we added in support for MPI parallelisation into hyperopt to enable it to search the space in parallel across all cores



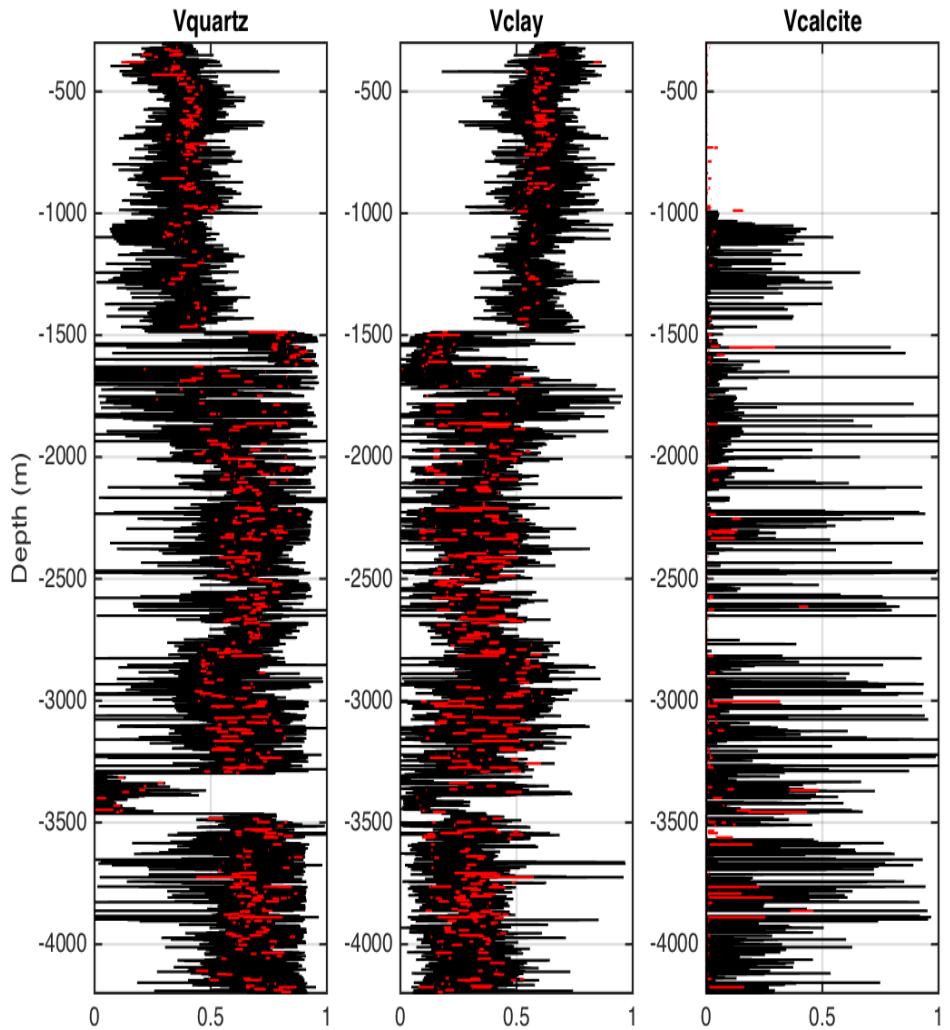
- Based on this parallelisation we can run around 2,500 parameter searches in a day now on a node of Cirrus
 - We can scale up to more nodes (Cirrus has 280 nodes, so based on this it would be possible to do around 20,000 searches in day) but this isn't needed at the moment
 - We tend to find the optimum parameters in around 500 searches, so in reality it means we only need to run the script for about 4 or 5 hours

Mineralogy: Test on a single well

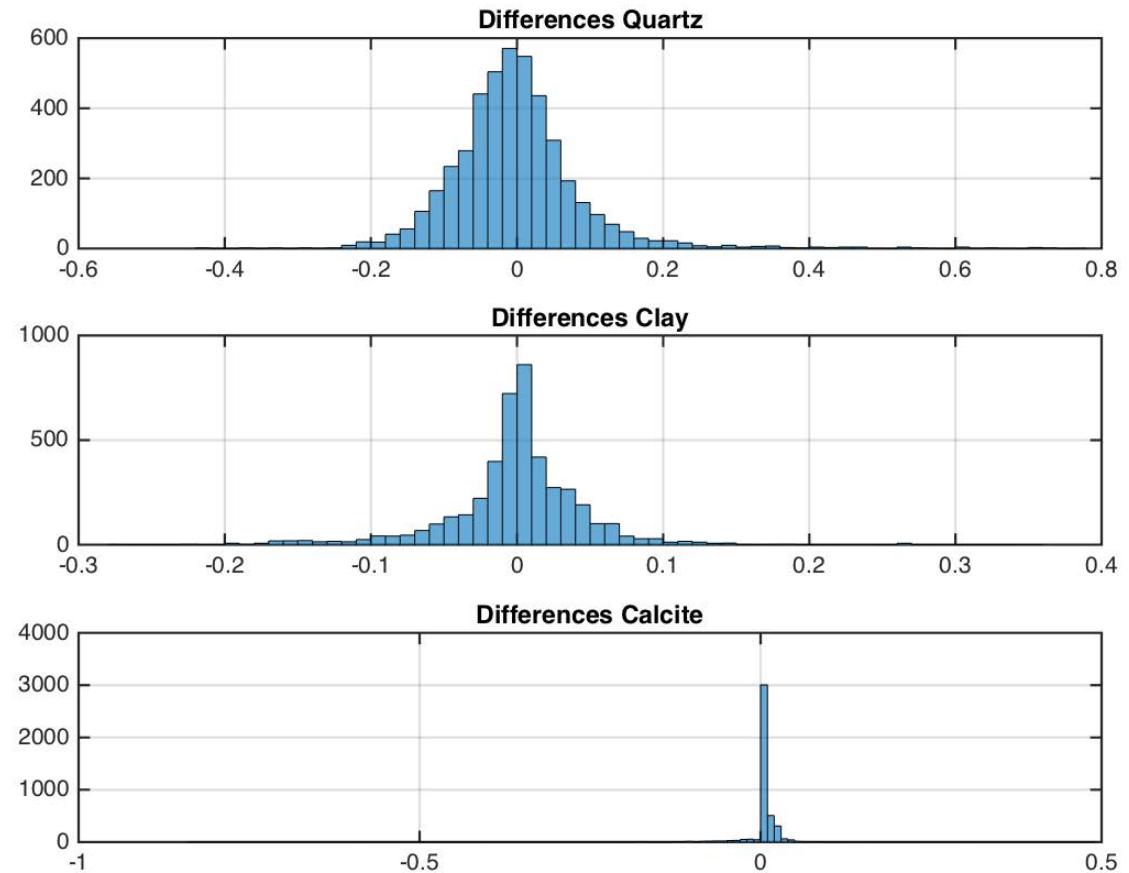


- Look at a single well: for this case the training data is as good as it can be !
- 80% of the points for training, 20% for testing.

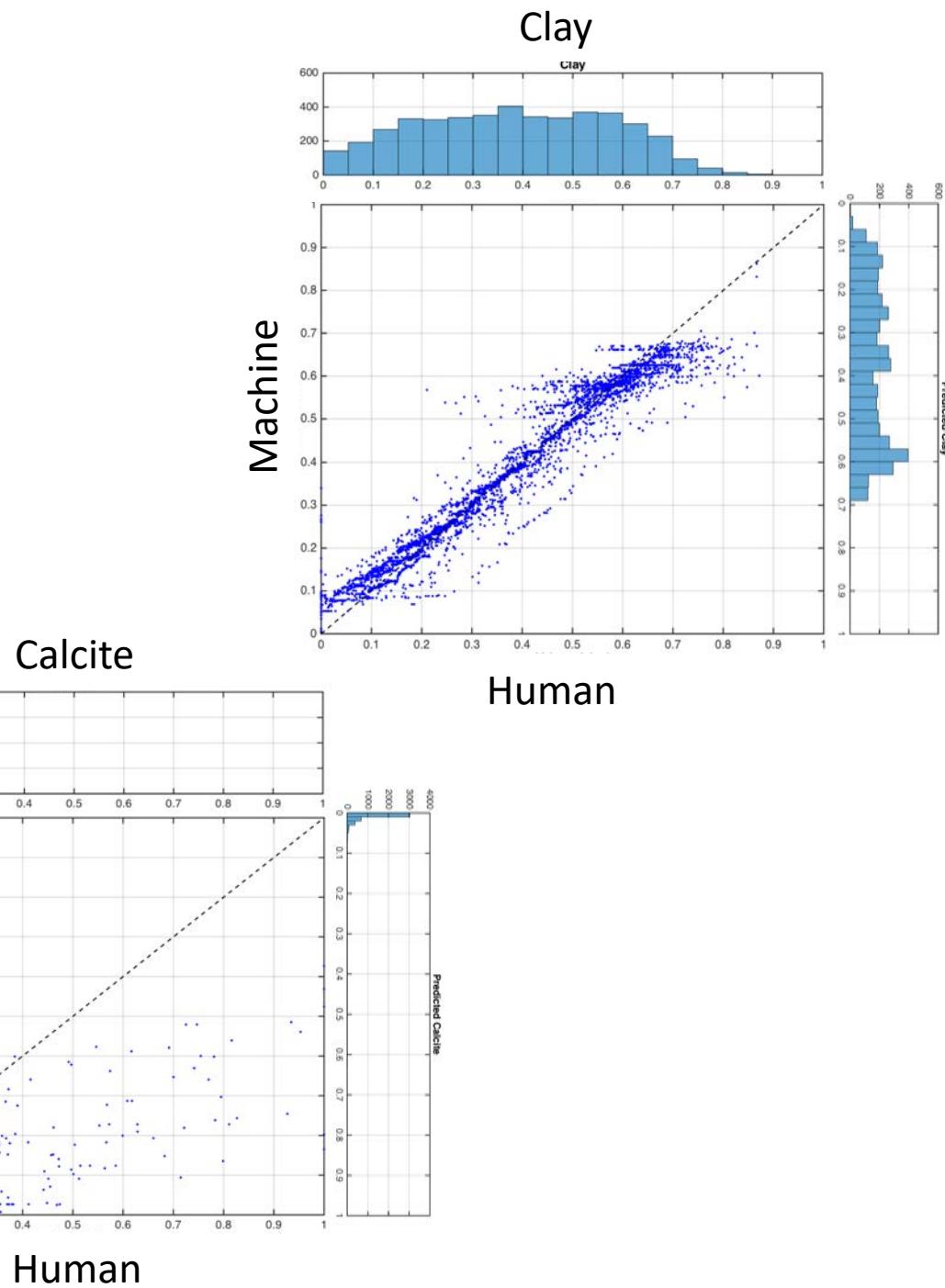
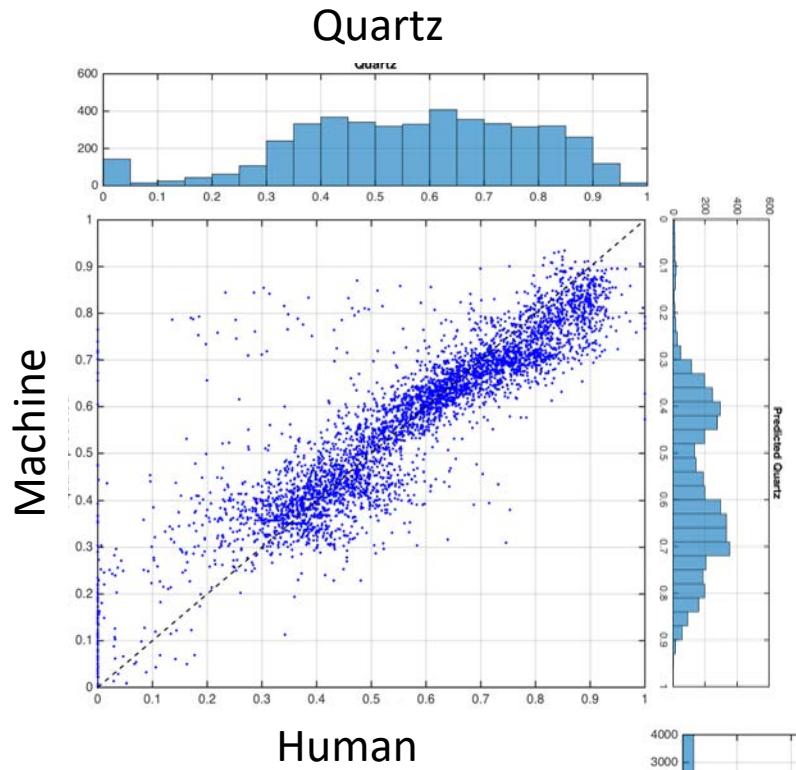
Results



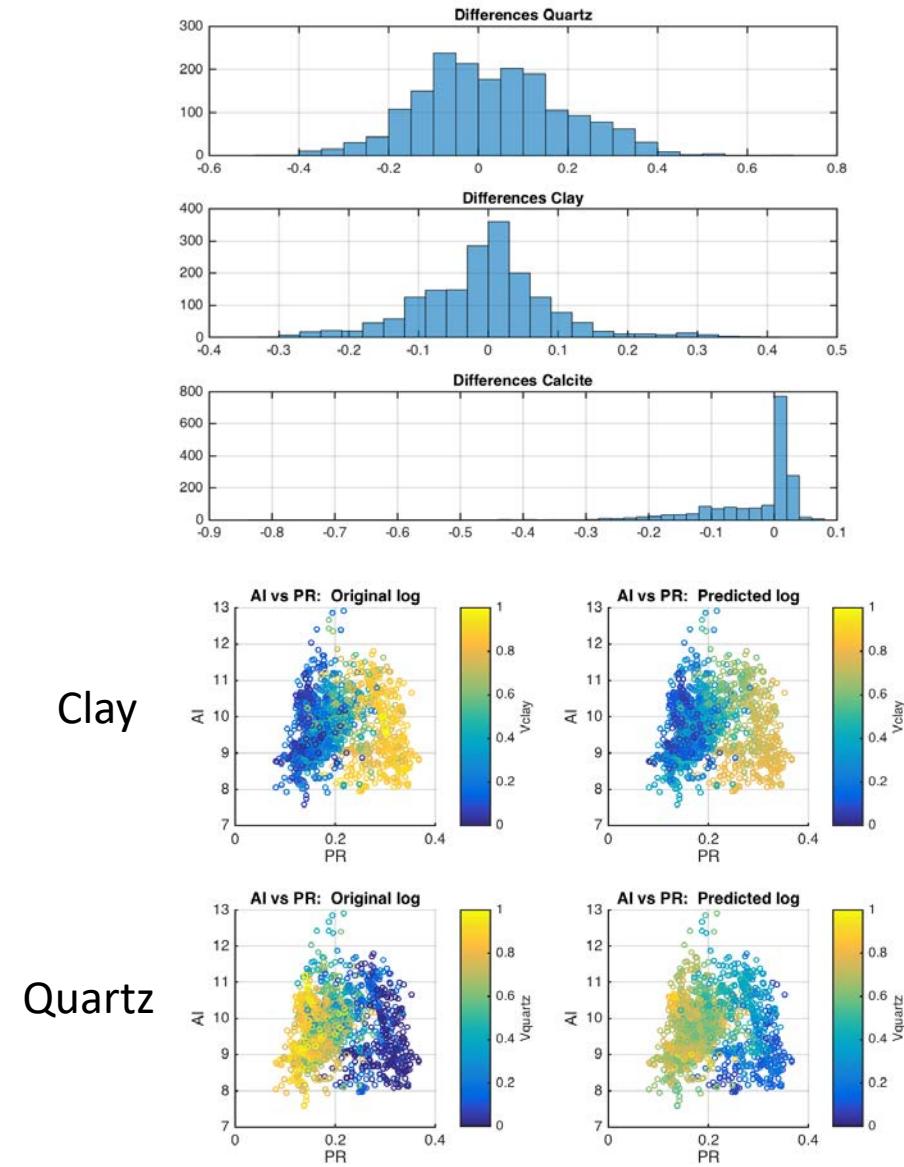
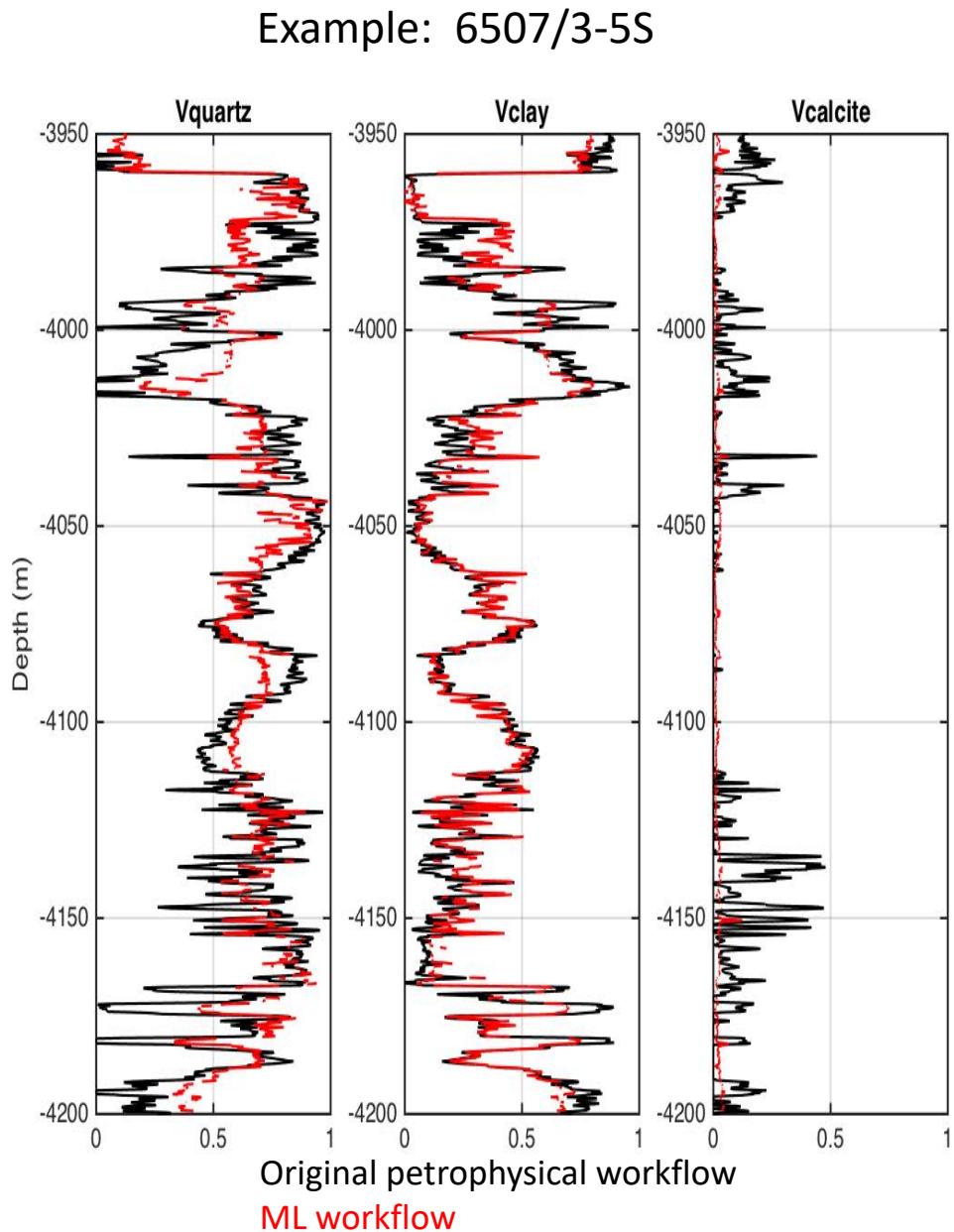
Original petrophysical workflow
ML workflow



Results

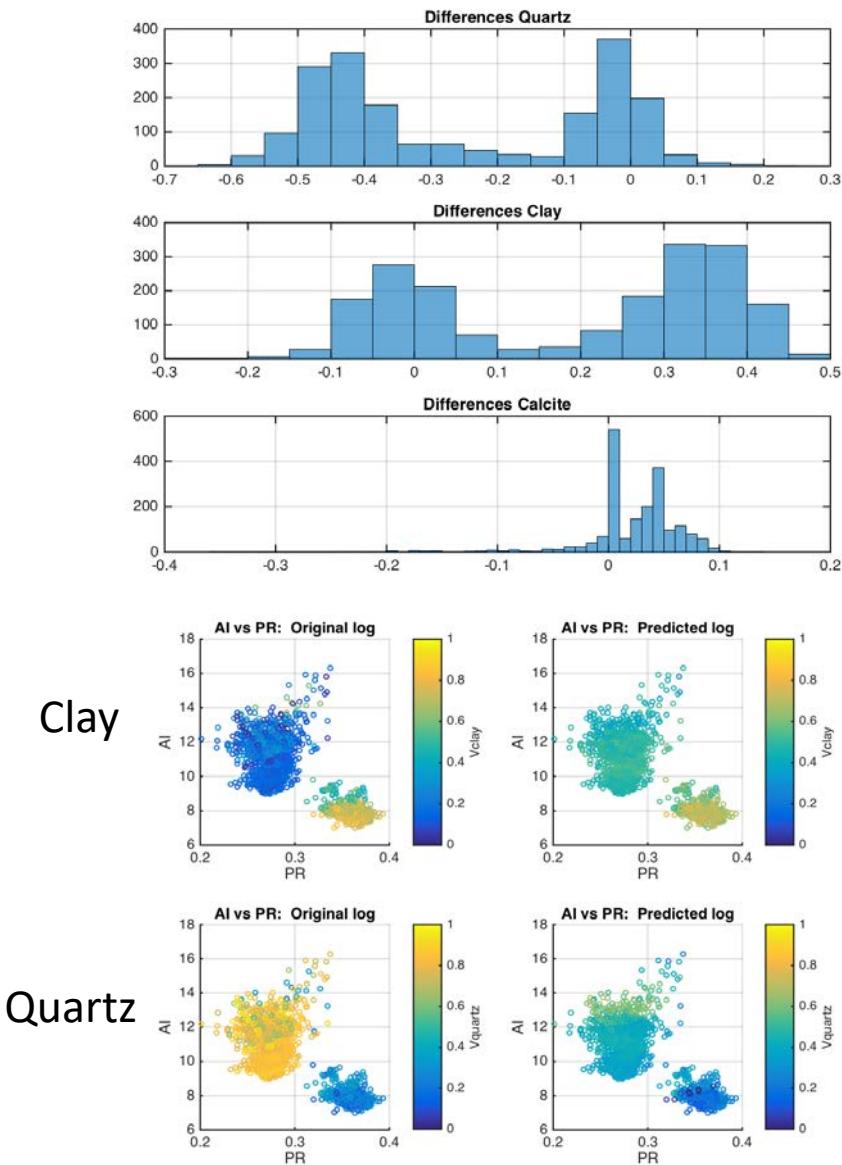
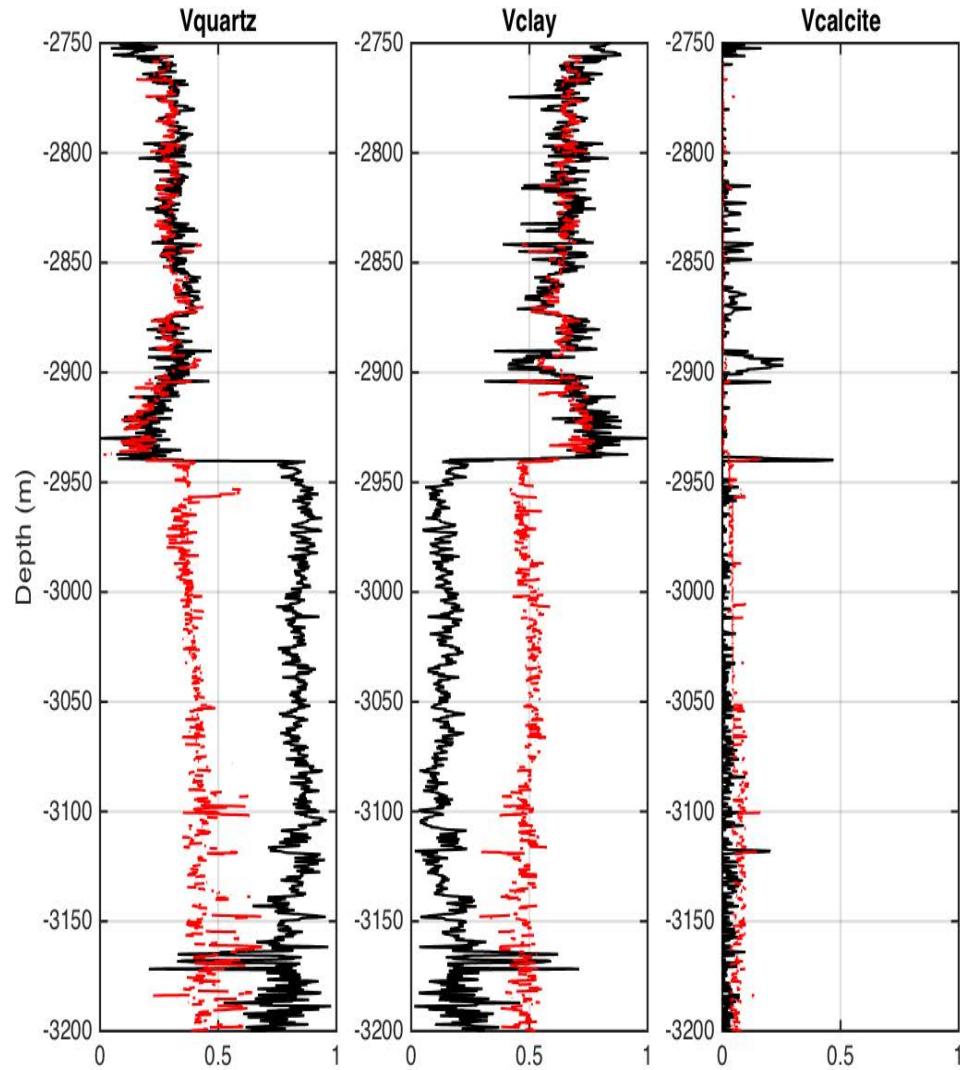


So far so good: what about predictions across wells ?

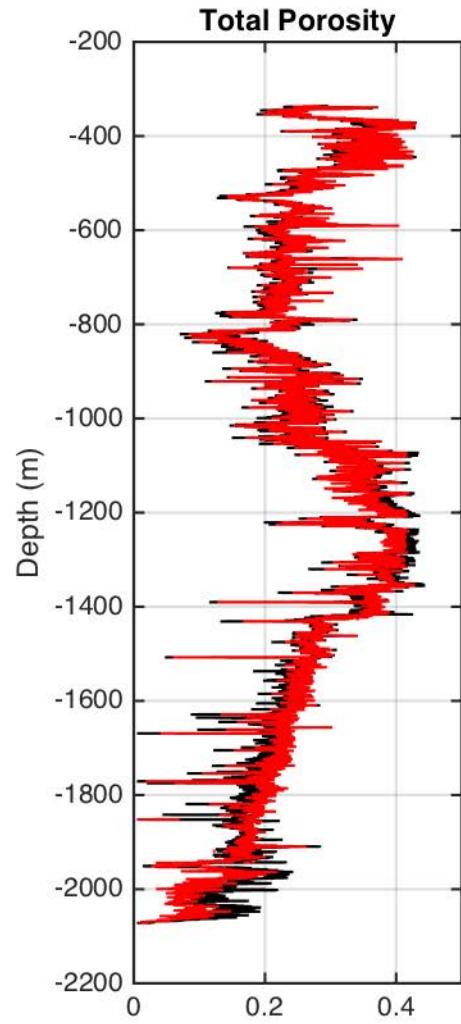


So far so good: what about predictions across wells ?

Example: 6306/5-2

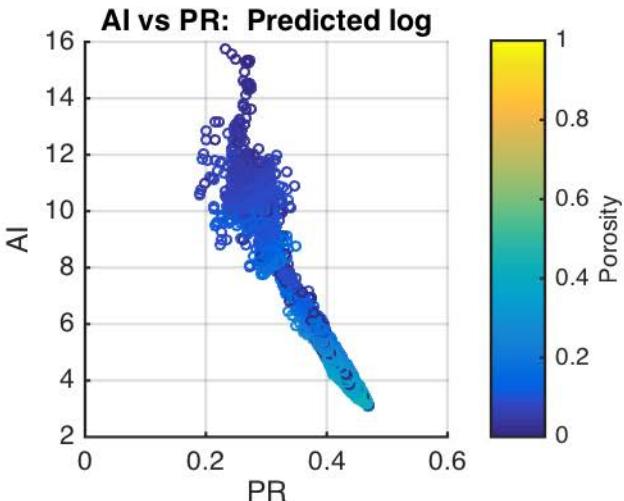
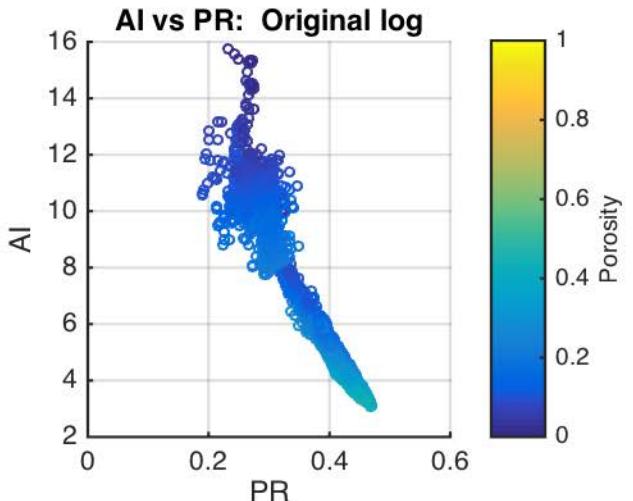
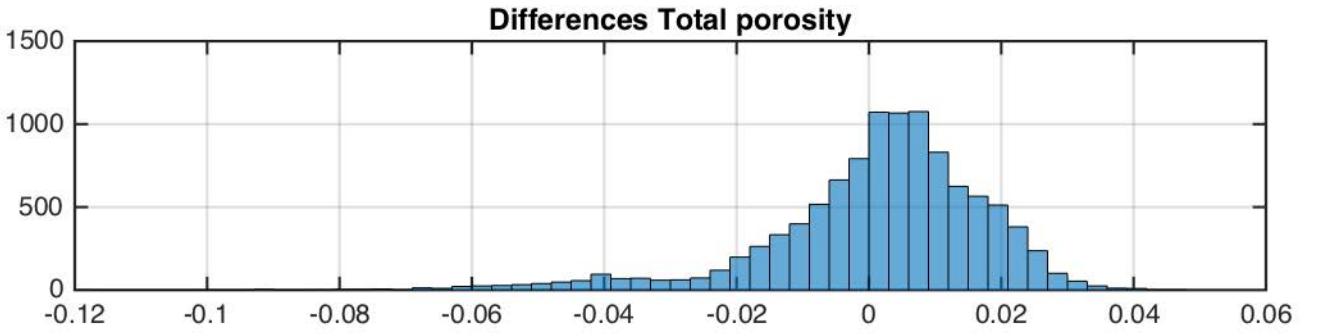
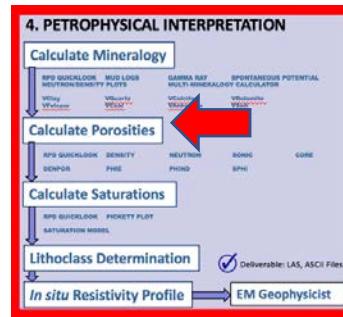


Predicting total porosity

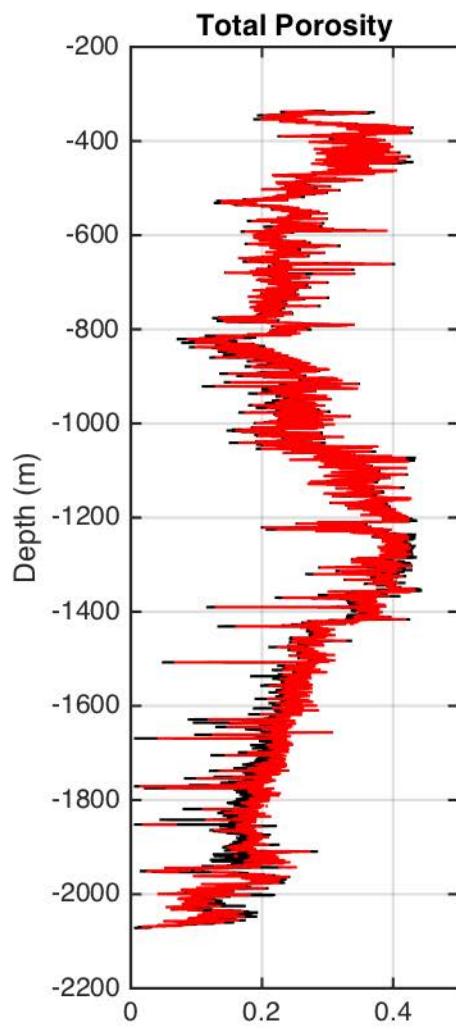


Example: 6306/6_2

Predictors: GR, Nphi, Resistivity, Rhob, Vp



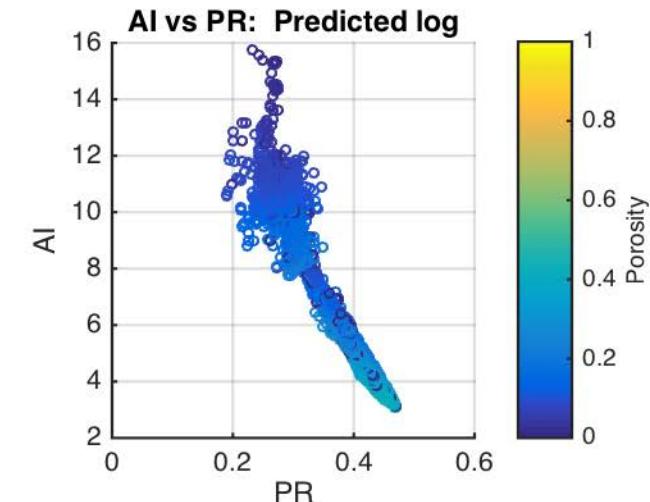
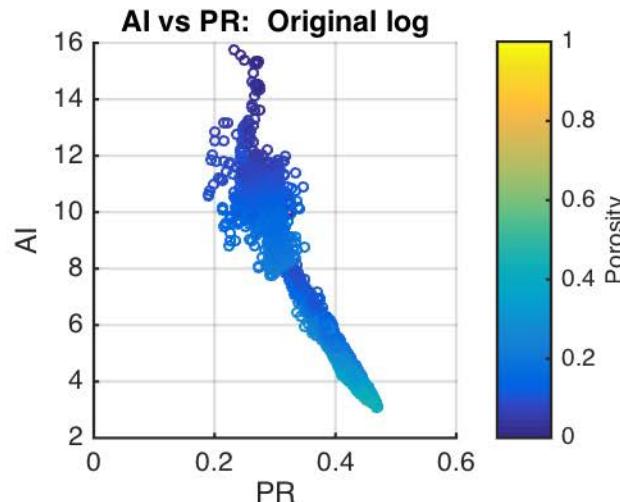
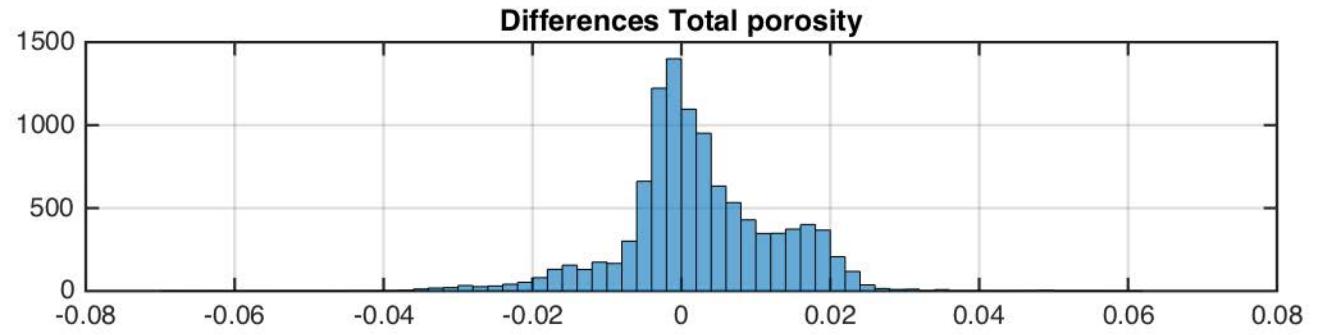
Predicting total porosity: add mineralogy



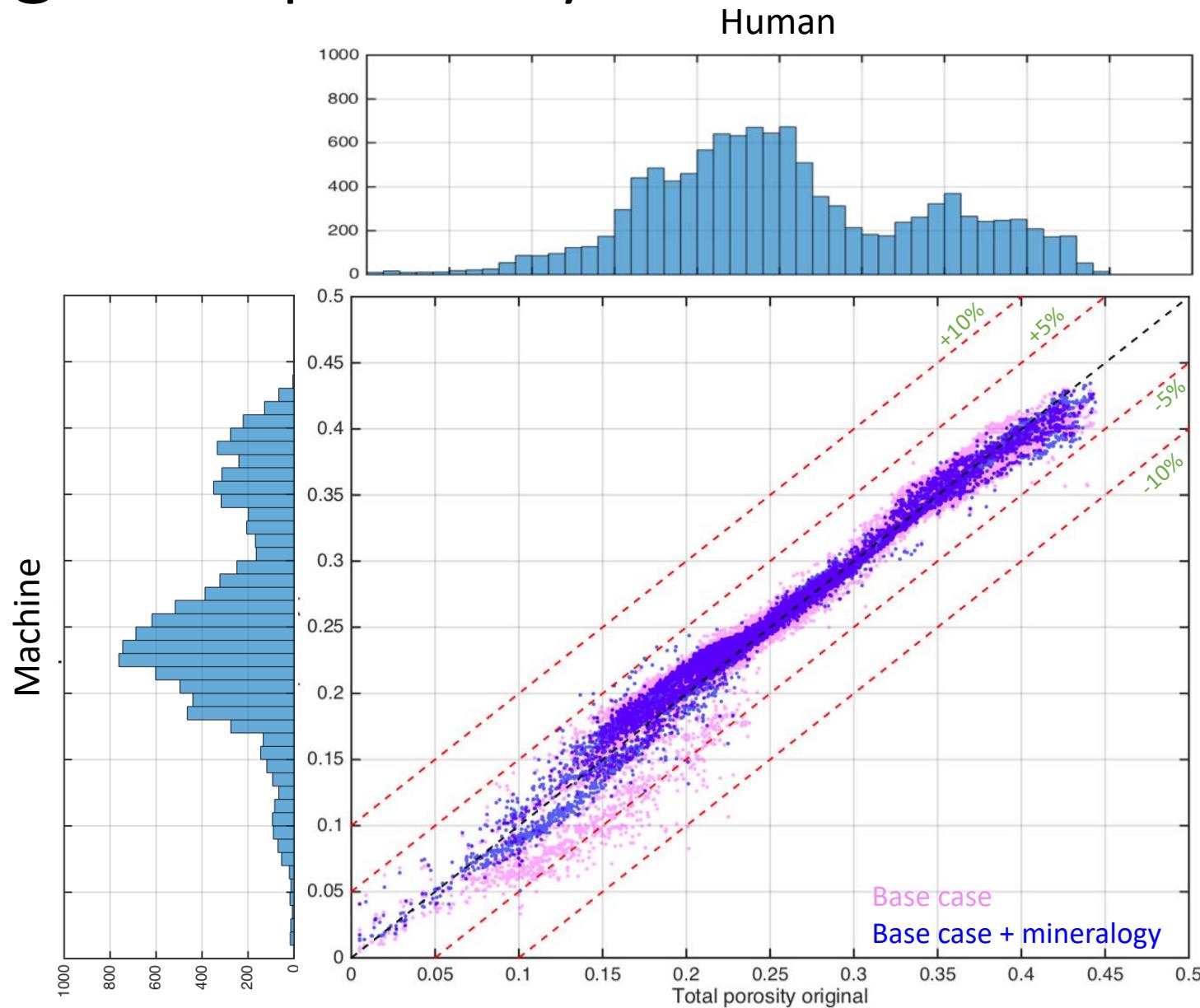
Original petrophysical workflow
ML workflow

Example: 6306/6_2

Predictors: GR, Nphi, Resistivity, Rhob, Vp, full mineral suite

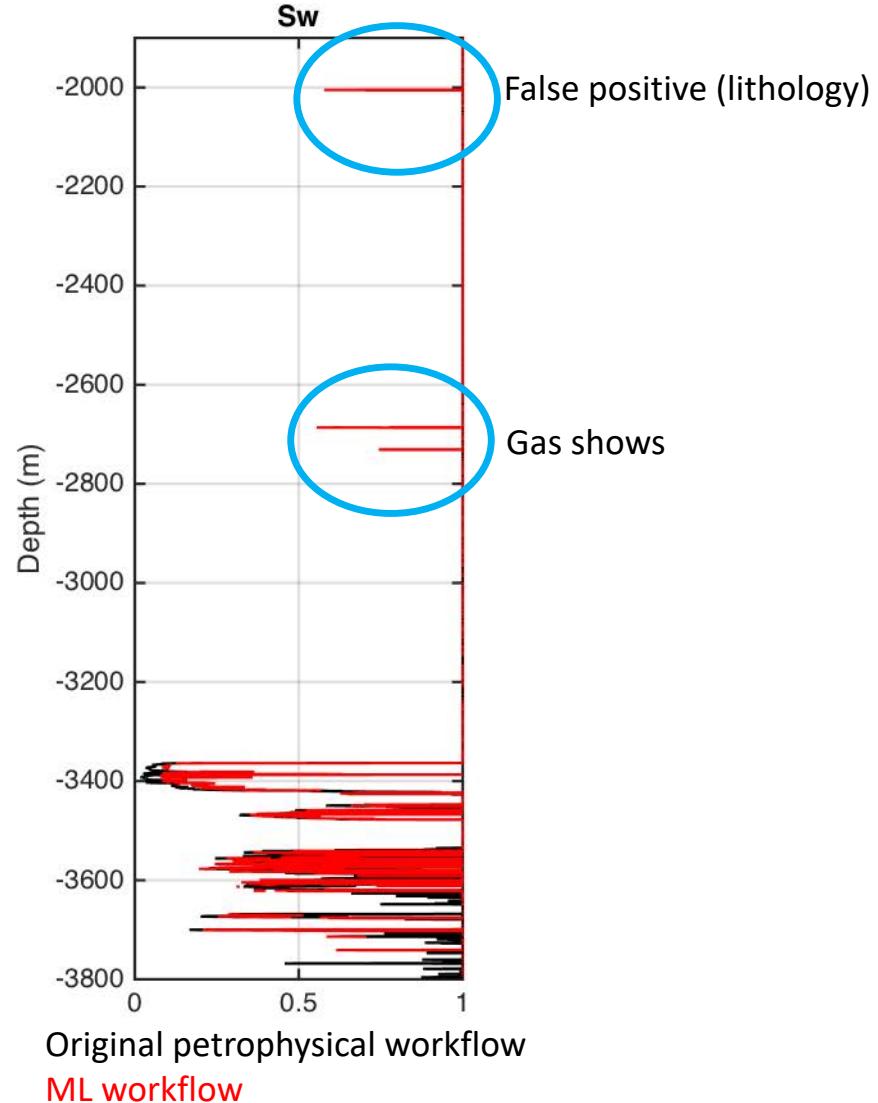


Predicting total porosity

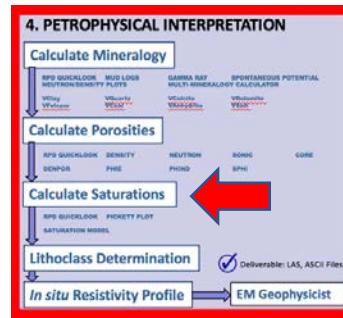
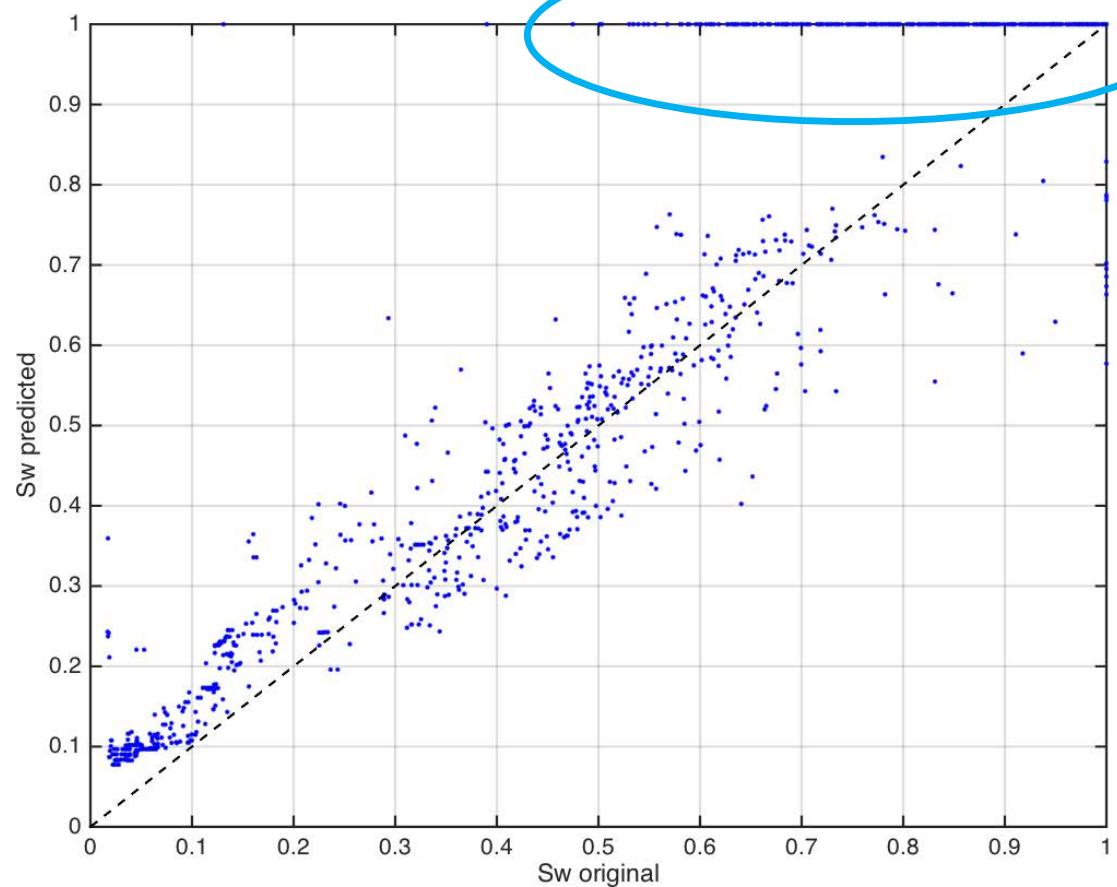


Predicting saturation: S_w

Well 6507/3_3



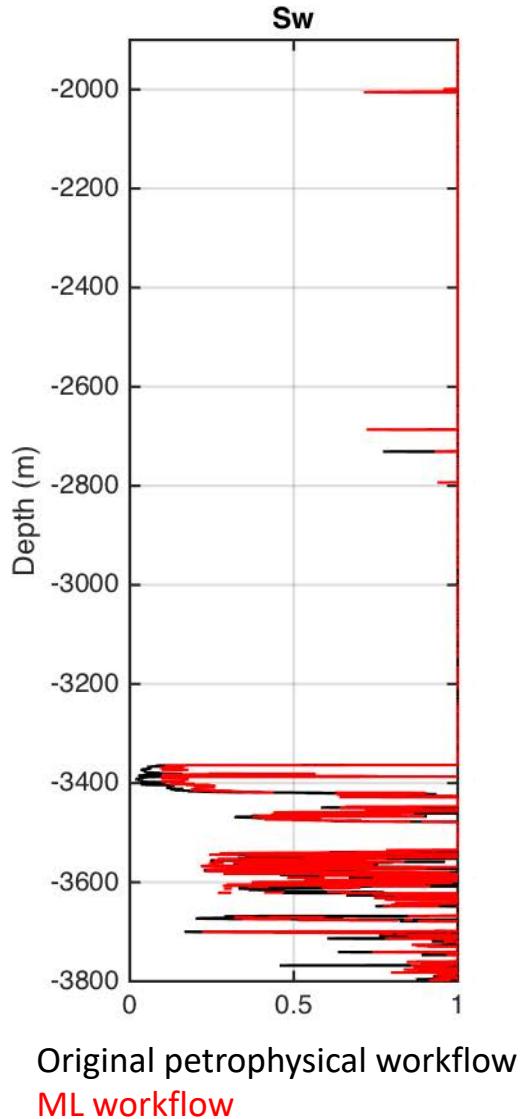
Training data bias: far more water than hydrocarbon in the training dataset



Balance the training data

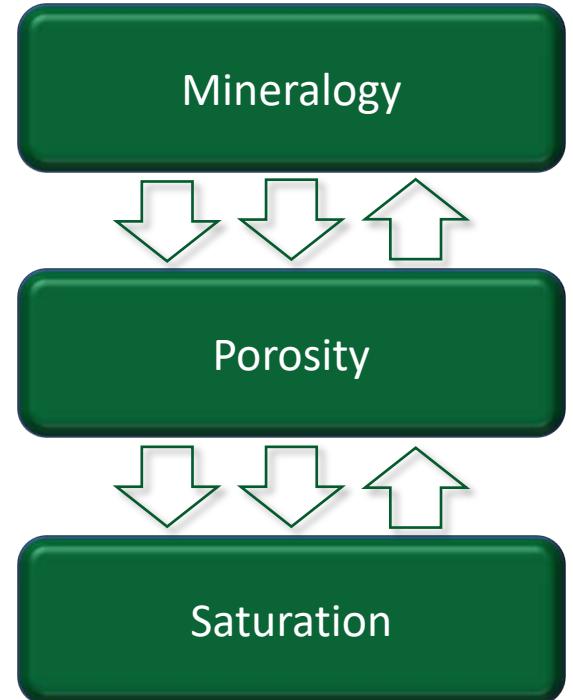
Well 6506/3_3

Predictors: GR, Nphi, Resistivity, Rhob, Vp, total porosity

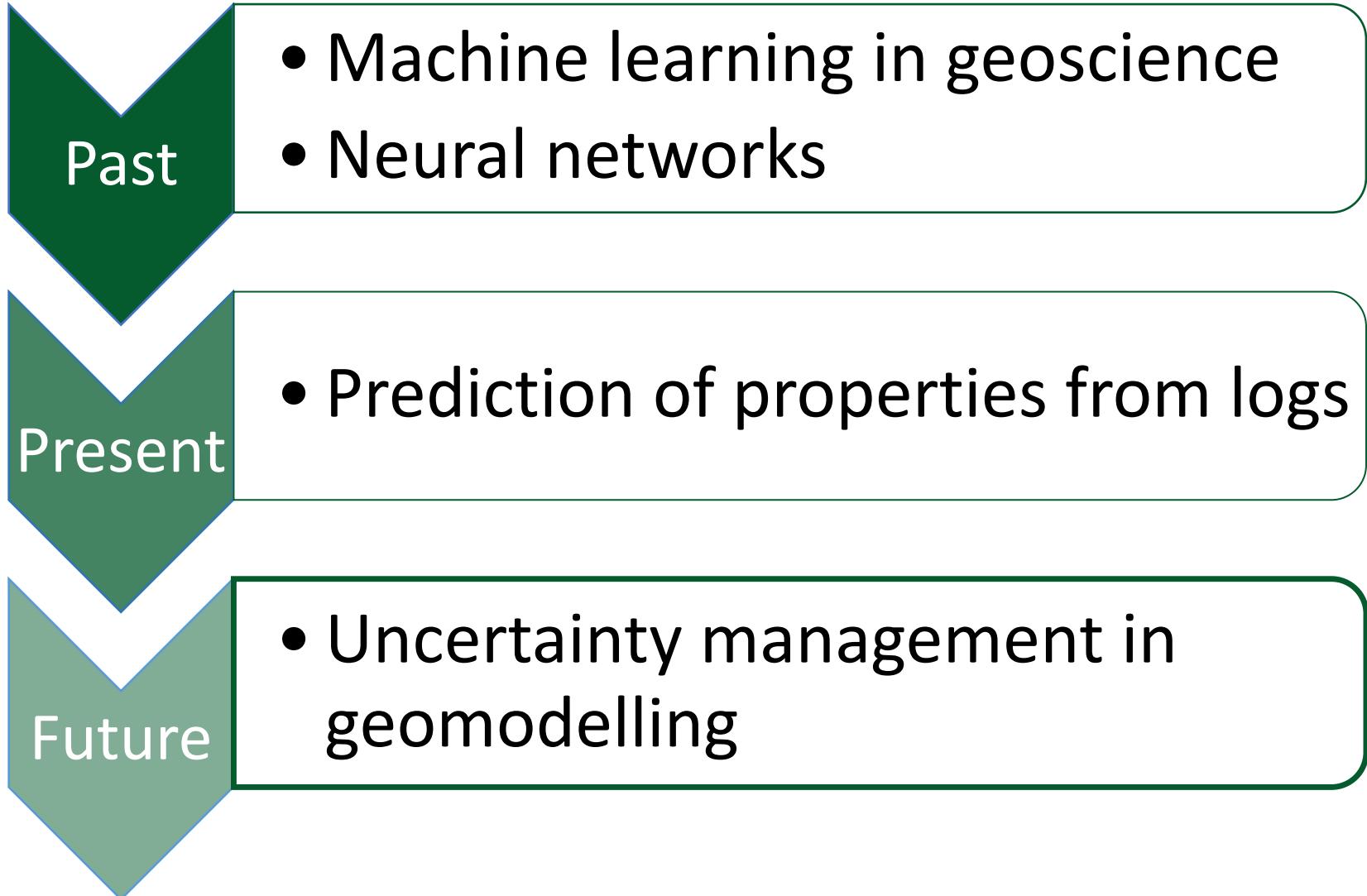


Observations

- Can we use machine learning to streamline workflows ?
 - Yes – results are promising
 - ...but we need to understand the effect of biased training data and assess how general a model can be built.
 - Needs to be iterative
- How good a model do we need ?
 - It doesn't need to be perfect to be useful. None of our models are perfect !
 - We're comparing against a human interpretation – who is right ?
 - How good is good enough and how do we characterize that ?



Overview



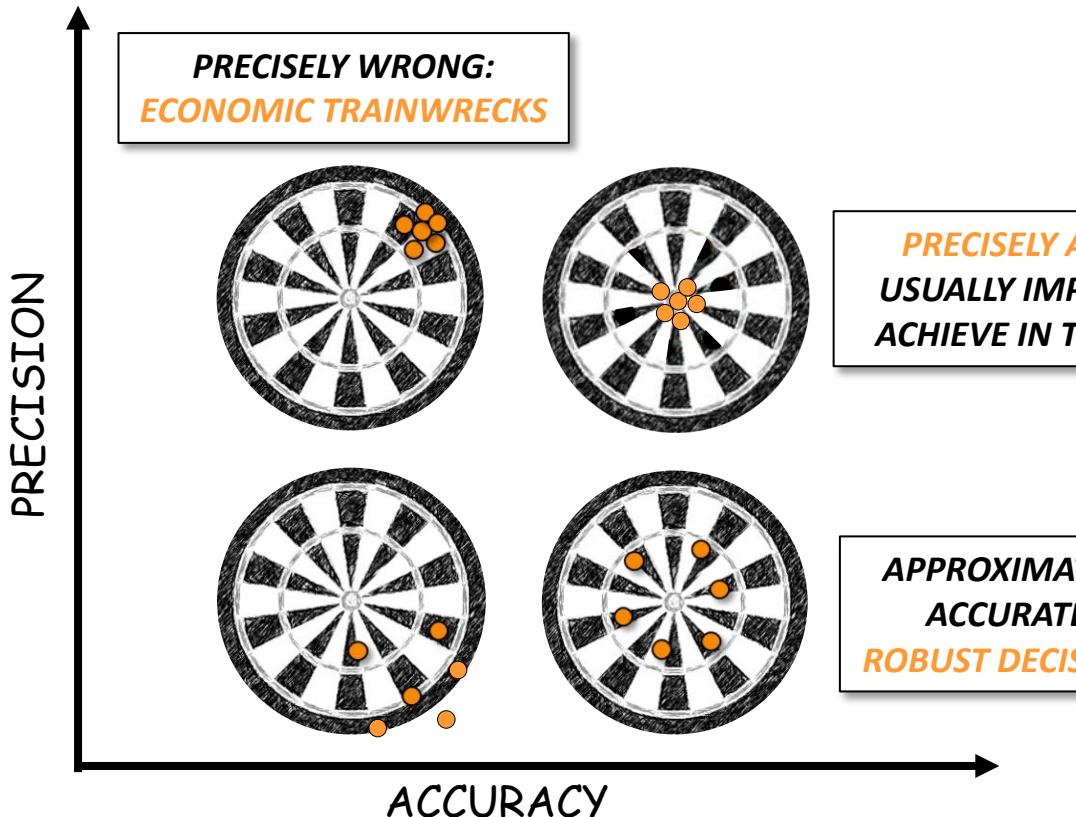
Geo-modelling in the oil and gas industry



*“all things equal, the best geologist
is the one who’s seen the most rocks”*

Geology is a science of case based reasoning, but **not** one where we can determine a **precise** answer

Accuracy versus precision

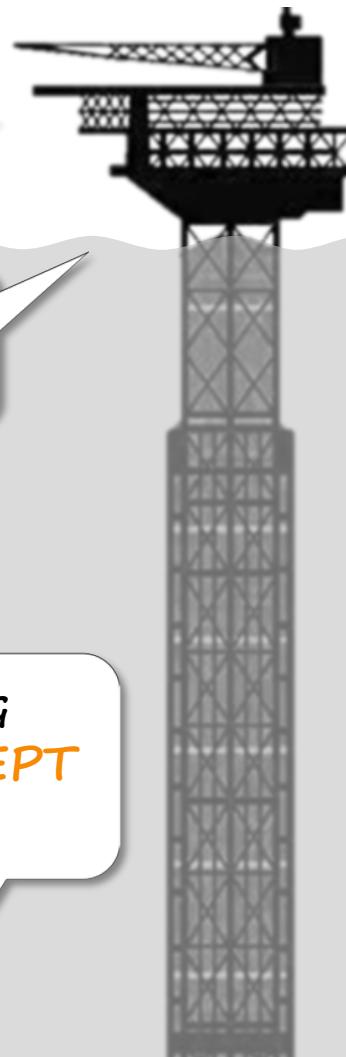


\$4 BILLION
DEVELOPMENT

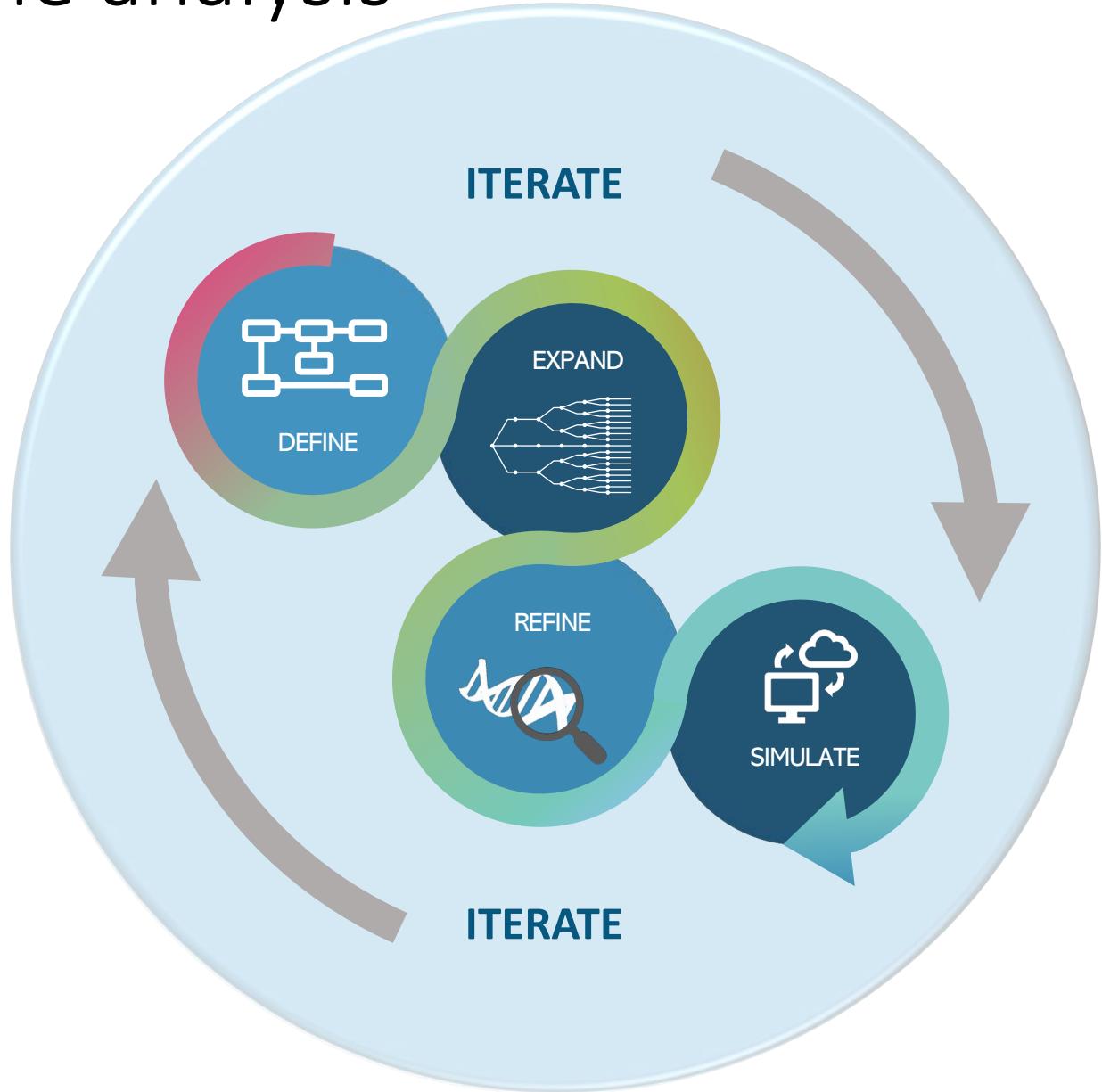
OPERATING AT
4% CAPACITY

BECAUSE UNDERLYING
GEOLOGICAL CONCEPT
WAS FLAWED

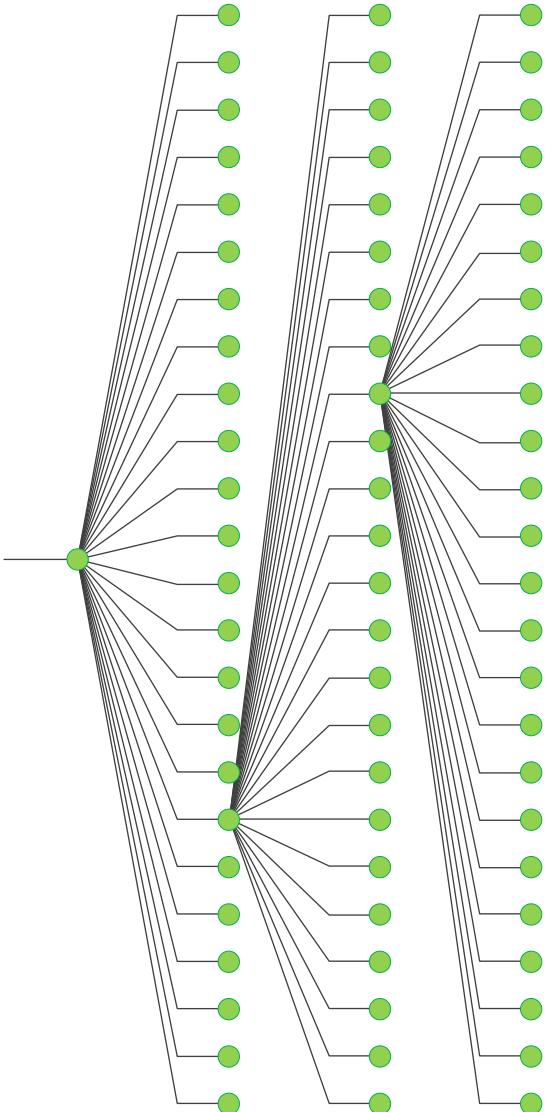
FAILURES ARE
EXPENSIVE



Model ensemble analysis



Creating the ensemble

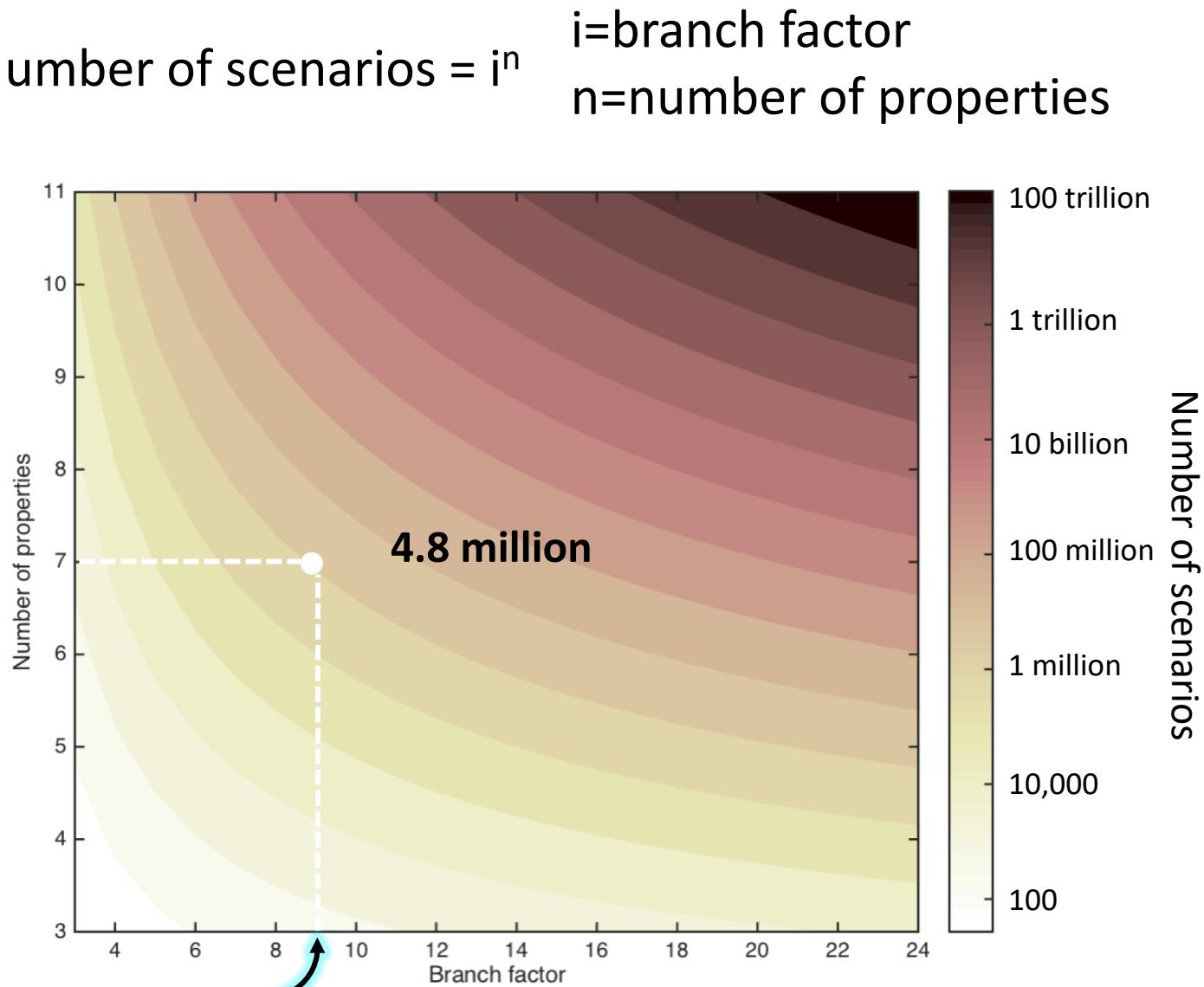


Number of scenarios = i^n

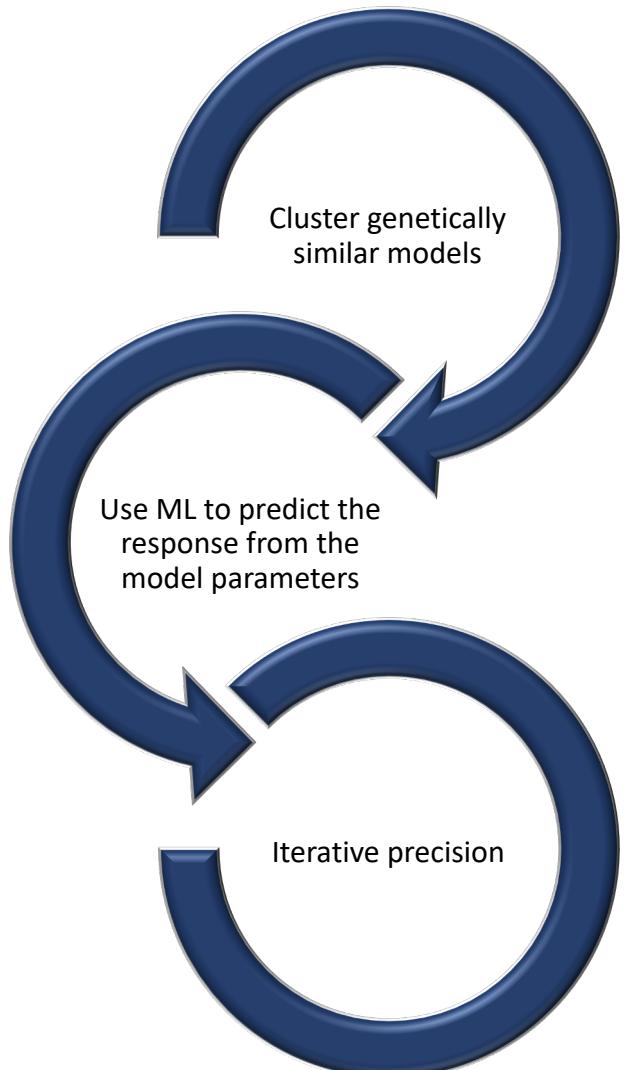
Grid
Facies
NTG
Porosity
XY Perm
Kv/Kh
Swirr



Percentiles (P10 to P90)



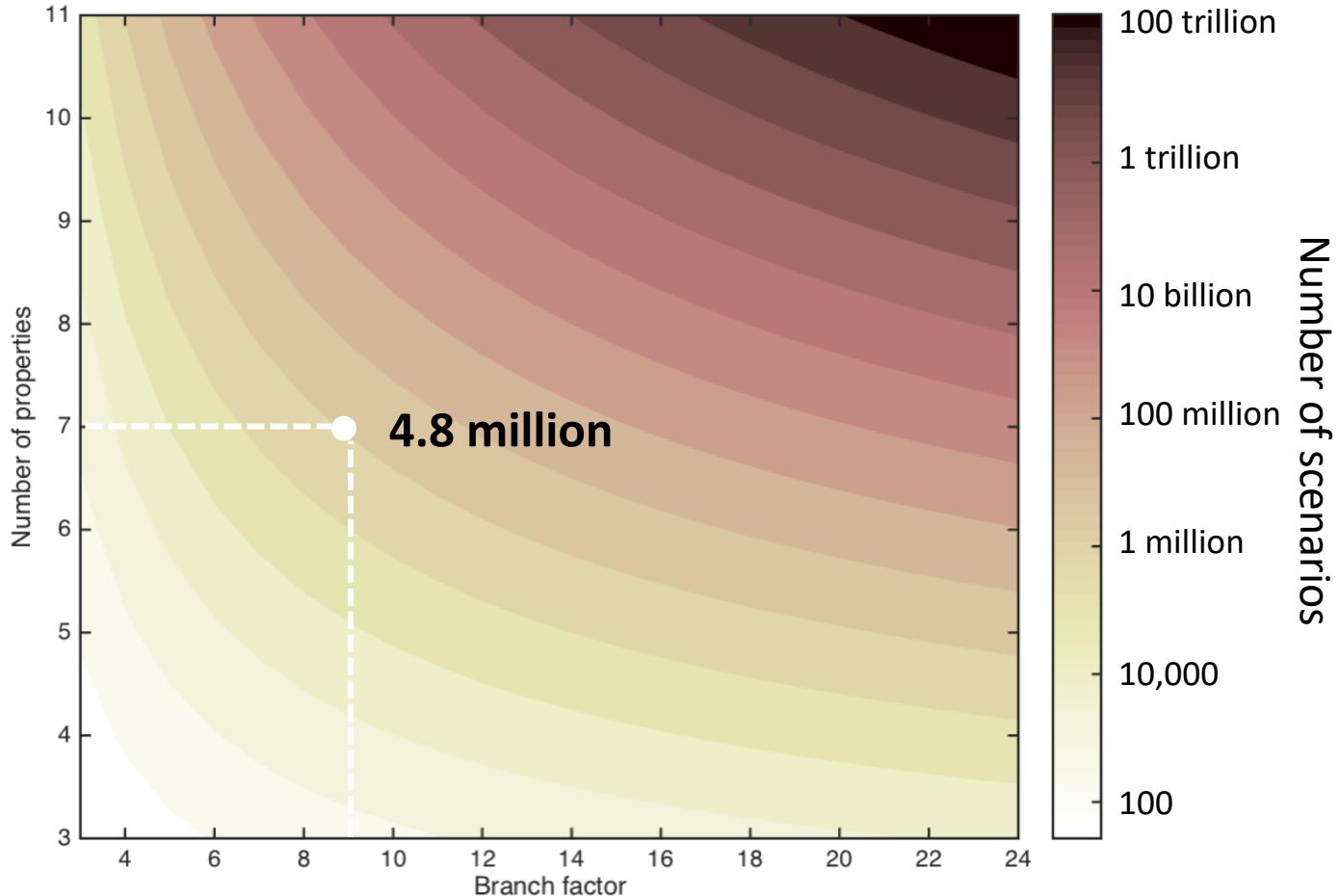
Managing the ensemble !



Number of scenarios = i^n

i =branch factor

n =number of properties



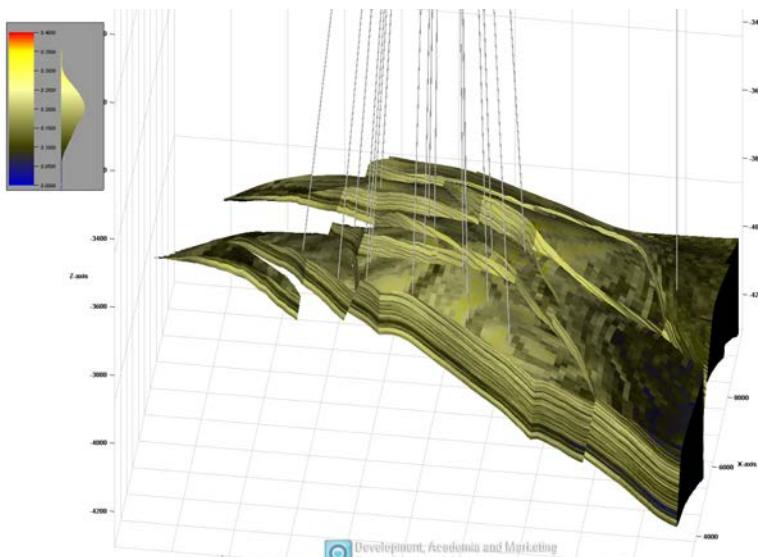
A really (really) simple example

Suppose we want to calculate the amount of oil in a reservoir:

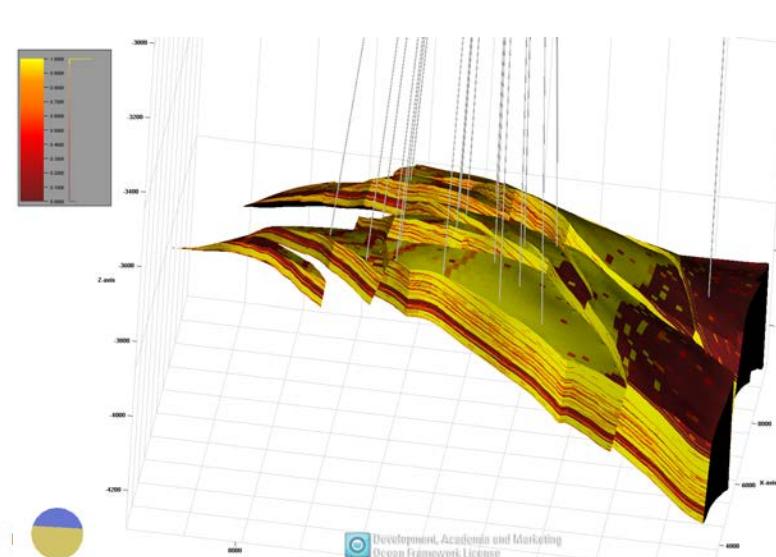
$$\text{Oil in place} = \text{Net to Gross} \times \text{Porosity} \times \text{Oil saturation}$$

Number of scenarios for each property: 24

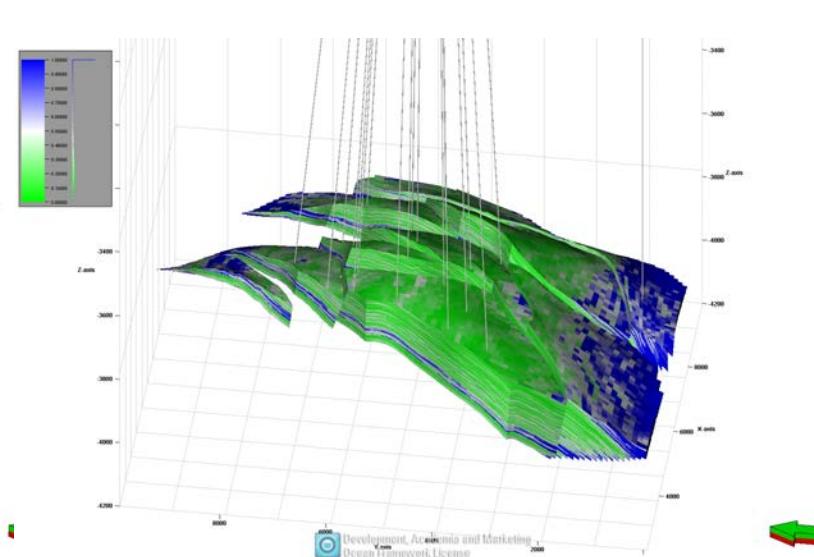
$$\text{Total number of oil in place values} = 24^3 = 13824$$



Porosity



NTG



Saturation

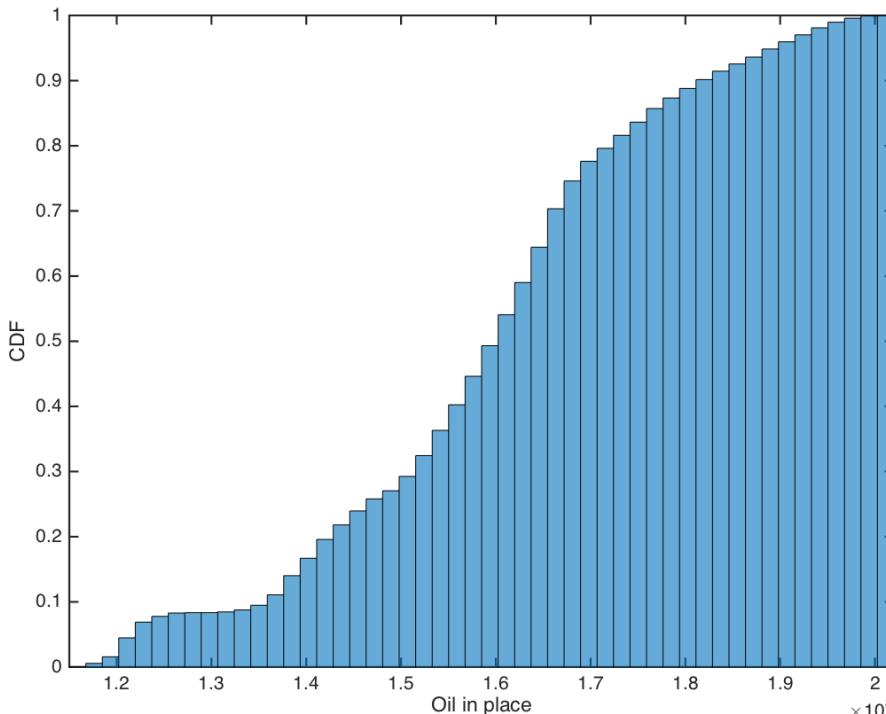
A really (really) simple example

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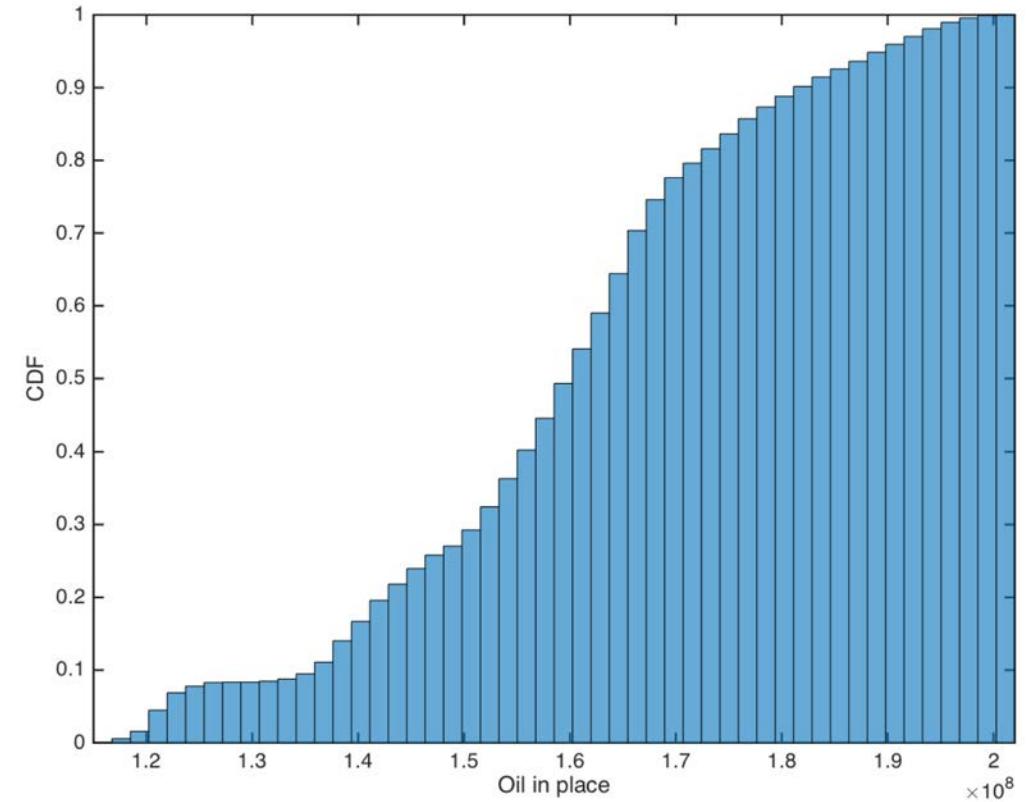
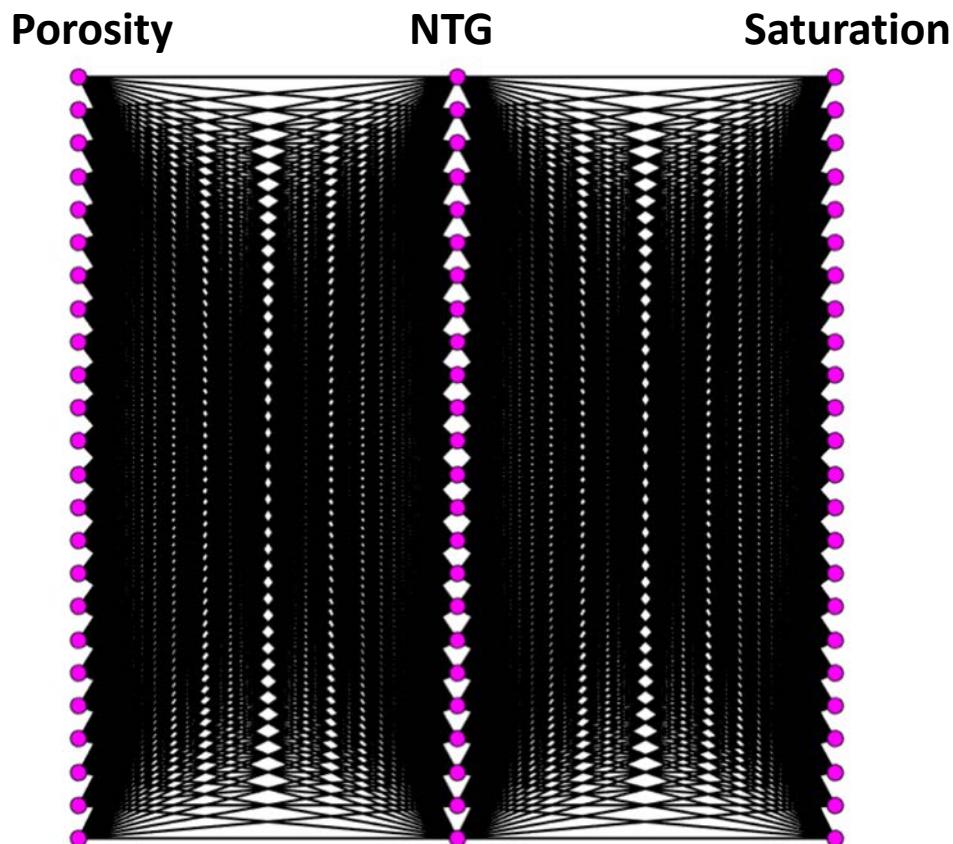
Number of scenarios for each property: 24

Total number of oil in place values = $24^3 = 13824$

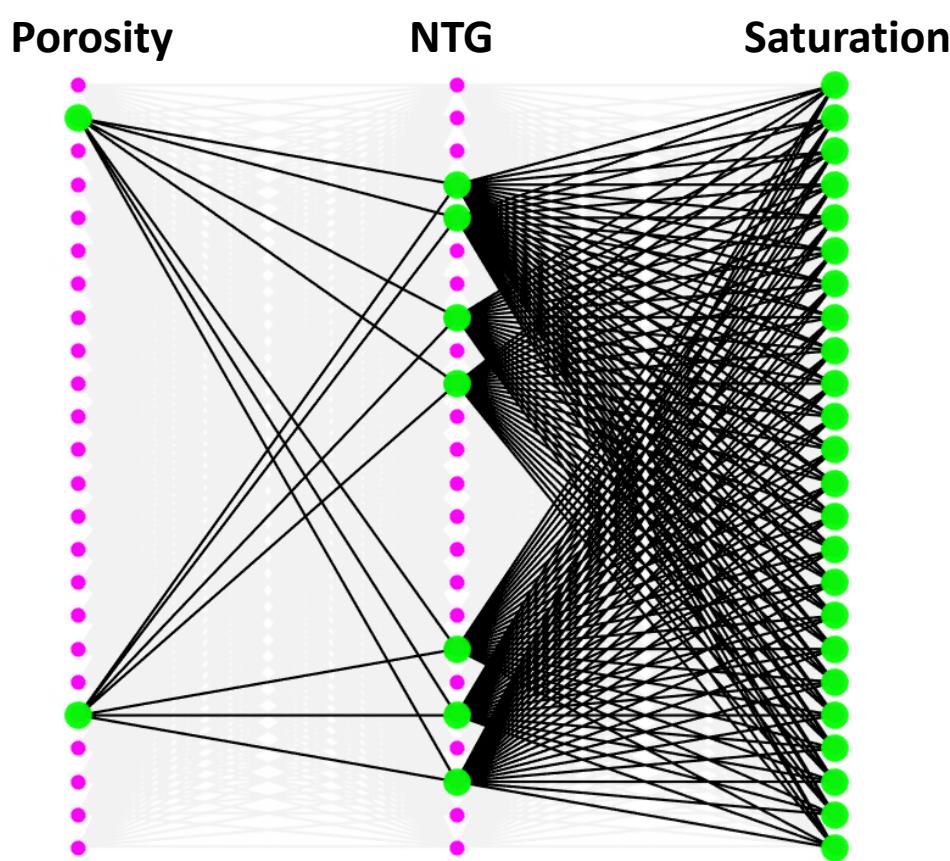


How many models do we need ?

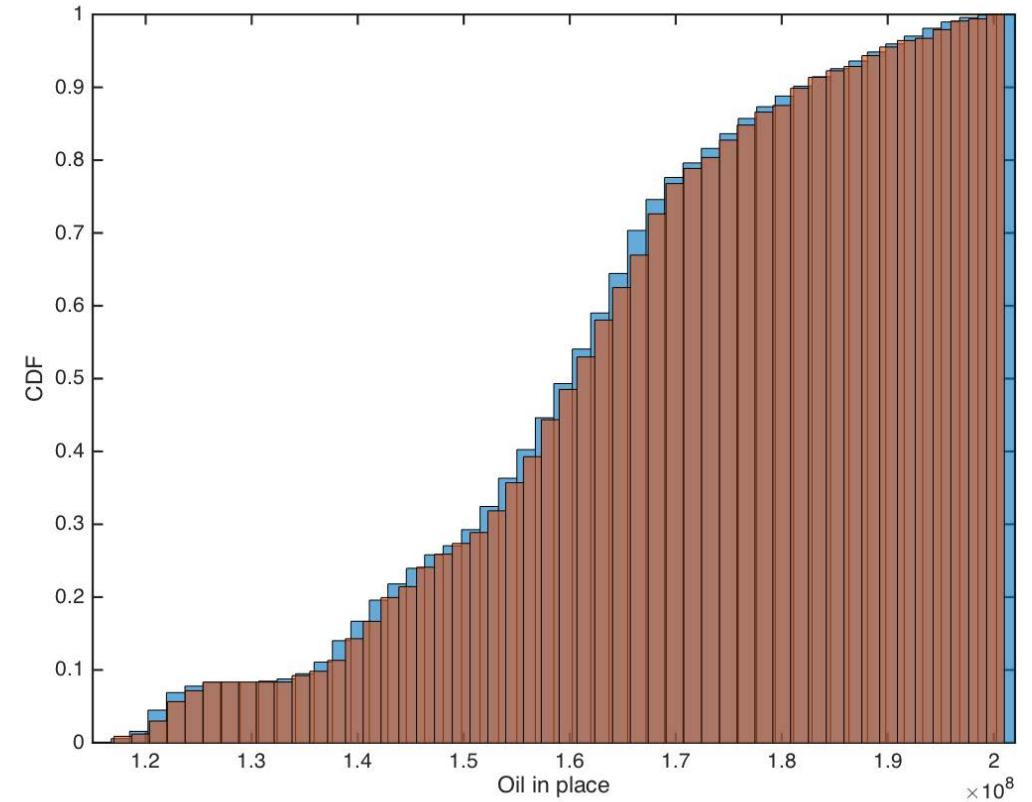
13824 models



How many models do we need ?



13824 models
366 models



Really simple example.....but shows that the approach is promising.

Questions and challenges

- How much training data is required? WHAT training data is needed ?
- How general will a trained machine learning approach be? For example, whereas you can imagine that results may be good within a single geological basin, will they be equally good on the other side of the planet?
- How accurate will a machine generated interpretation be? It does not need to be perfect to be useful: something that is right 70-80% of the time could still represent a huge time saving in the workflow.
- Can we characterize/understand the uncertainty in the results?