
Segmenting neutron images using machine learning

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**CHAPTER
ONE**

LECTURE OUTLINE

In this lecture about machine learning to segment neutron images we will cover the following topics

1. Introduction
2. Limited data problem
3. Unsupervised segmentation
4. Supervised segmentation
5. Final problem: Segmenting root networks using convolutional NNs
6. Future Machine learning challenges in NI

1.1 Getting started

If you want to run the notebook on your own computer, you'll need to perform the following step:

- You will need to install Anaconda
- Clone the lecture repository (in the location you'd like to have it)

```
git clone https://github.com/ImagingLectures/MLSegmentation4NI.git
```

- Enter the folder 'MLSegmentation'
- Create an environment for the notebook

```
conda env create -f environment.yml -n MLSeg4NI
```

- Enter the environment

```
conda env activate MLSeg4NI
```

- Start jupyter and open the notebook `lecture/ML4NeutronImageSegmentation.ipynb`
- Use the notebook
- Leave the environment

```
conda env deactivate
```

1.2 Importing needed modules

This lecture needs some modules to run. We import all of them here.

```
import matplotlib.pyplot as plt
import seaborn as sn
import numpy as np
import pandas as pd
import skimage.filters as flt
import skimage.io as io
import matplotlib as mpl

from sklearn.cluster import KMeans
from sklearn.neighbors import KNeighborsClassifier
from sklearn.metrics import confusion_matrix
from sklearn.datasets import make_blobs

from matplotlib.colors import ListedColormap
from matplotlib.patches import Ellipse
from lecturesupport import plotsupport as ps

import scipy.stats as stats
import astropy.io.fits as fits

import keras.metrics as metrics
import keras.losses as loss
import keras.optimizers as opt
from keras.models import Model
from keras.layers import Input, Conv2D, MaxPooling2D, UpSampling2D, concatenate

%matplotlib inline

from IPython.display import set_matplotlib_formats
set_matplotlib_formats('svg', 'png')
# plt.style.use('seaborn')
mpl.rcParams['figure.dpi'] = 300
```

Using TensorFlow backend.

```
import importlib
importlib.reload(ps);
```

**CHAPTER
TWO**

INTRODUCTION

In this introduction part we give a starting point for the lecture. With topics like.

- Introduction to neutron imaging
 - Some words about the method
 - Contrasts
- Introduction to segmentation
 - What is segmentation
 - Noise and SNR
- Problematic segmentation tasks
 - Intro
 - Segmentation problems in neutron imaging

From these topics we will go into looking at how different machine learning techniques can be used to segment images. An in particular images obtained in neutron imaging experiments

2.1 What is an image?

In this lecture we are going to analyze the contents of images, but first lets define what an image is.

A very abstract definition:

- **A pairing between spatial information (position)**
- **and some other kind of information (value).**

In most cases this is a two- or three-dimensional position (x,y,z coordinates) and a numeric value (intensity)

In reality it can be many things like a picture, a radiograph, a matrix presentation of a raster scan and so on. In the case of volume images we are talking about volume representations like a tomography or time series like movies.

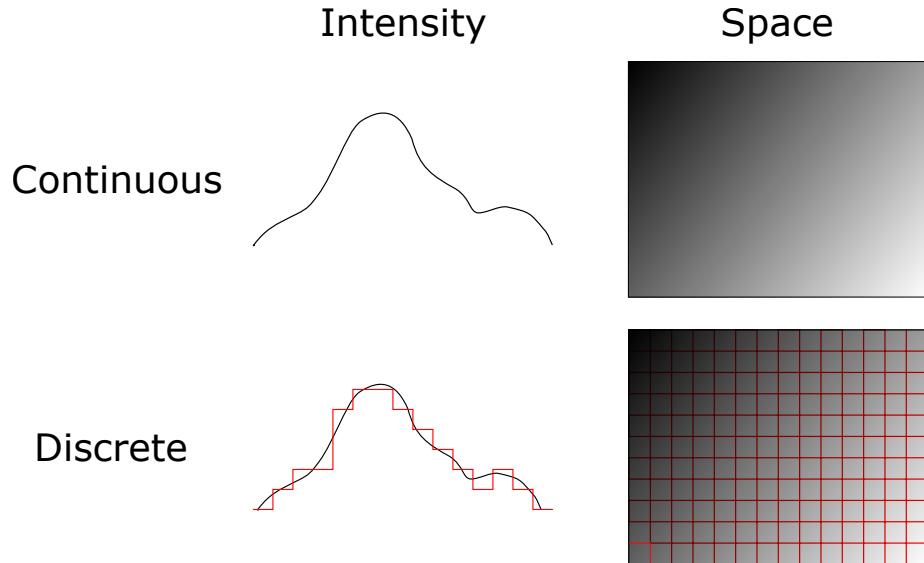


Fig. 2.1: The real world is sampled into discrete images with limited extent.

2.2 Science and Imaging

Images are great for qualitative analyses since our brains can quickly interpret them without large *programming* investments.

This is also the reason why image processing can be very frustrating. Seeing a feature in the image may be obvious to the human eye, but it can be very hard to write an algorithm that makes the same conclusion. In particular, when you are using traditional image processing techniques. This is a classic area for machine learning, these methods are designed and trained to recognize patterns and features in the images. Still, they can only perform well on the type of data they are trained with.

2.2.1 Proper processing and quantitative analysis is however much more difficult with images.

- If you measure a temperature, quantitative analysis is easy, $T = 50K$.
- If you measure an image it is much more difficult and much more prone to mistakes,
 - subtle setup variations may break your analysis process,
 - and confusing analyses due to unclear problem definition

2.2.2 Furthermore in image processing there is a plethora of tools available

- Thousands of algorithms available
- Thousands of tools
- Many images require multi-step processing
- Experimenting is time-consuming

2.3 Some word about neutron imaging

Neutron imaging is a method based on transmission of the neutron radiation through a sample, i.e. the fundamental information is a radiograph. In this sense it is very much similar to the more known x-ray imaging. The basic outline in the figure below show the components source, collimator, sample, and detector. The neutron beam is slightly divergent, but it can mostly be approximated by a parallel beam.

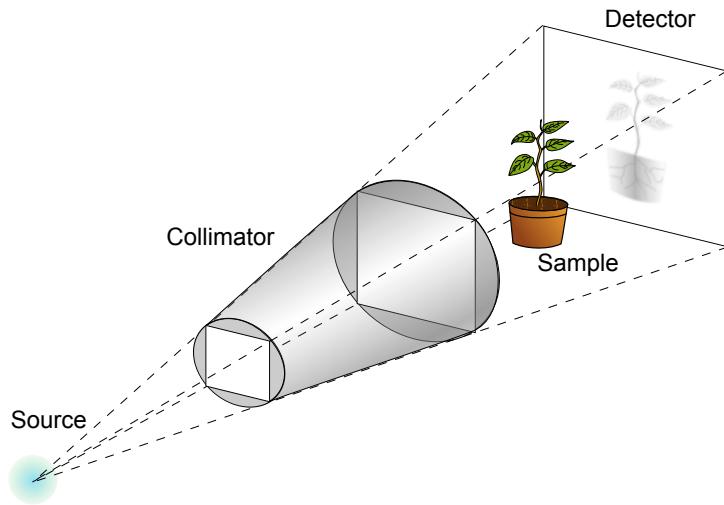


Fig. 2.2: Schematic transmission imaging setup.

The intensity in a radiographic image proportional to the amount of radiation that remains after it was transmitted through the sample.

The transmitted radiation is described by Beer-Lambert's law which in its basic form looks like

$$I = I_0 \cdot e^{-\int_L \mu(x)dx}$$

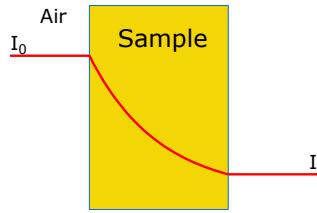


Fig. 2.3: Transmission plot through a sample.

Where $\mu(x)$ is the attenuation coefficient of the sample at position x . This equation is a simplification as no source has a uniform spectrum and the attenuation coefficient depends on the radiation energy. We will touch the energy dispersive property in one of the segmentation examples later.

Single radiographs are relatively rare. In most experiments the radiographs are part of a time series acquisition or projections for the 3D tomography reconstruction. In this lecture we are not going very much into the details about the imaging method as such. This is a topic for other schools that are offered.

2.3.1 Image types obtained with neutron imaging

| Fundamental information | Additional dimensions | Derived information |
|-------------------------|-----------------------|---------------------|
| 2D Radiography | Time series | q-values |
| 3D Tomography | Spectra | strain |
| | | Crystal orientation |

2.4 Neutron imaging contrast

The integral of Beer-Lambert's law turns into a sum for simple material compositions like in the sample in the figure below

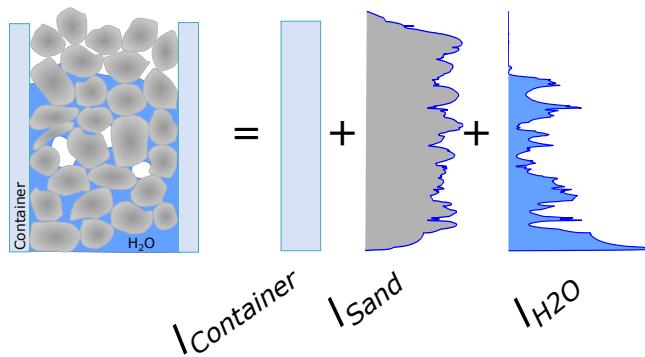


Fig. 2.4: The transmission through a sample can be divided into contributions for each material.

The amount of radiation passing through the sample depends on the attenuation coefficients which are material and radiation specific. The periodic systems below illustrate how different elements attenuate for neutrons and X-rays

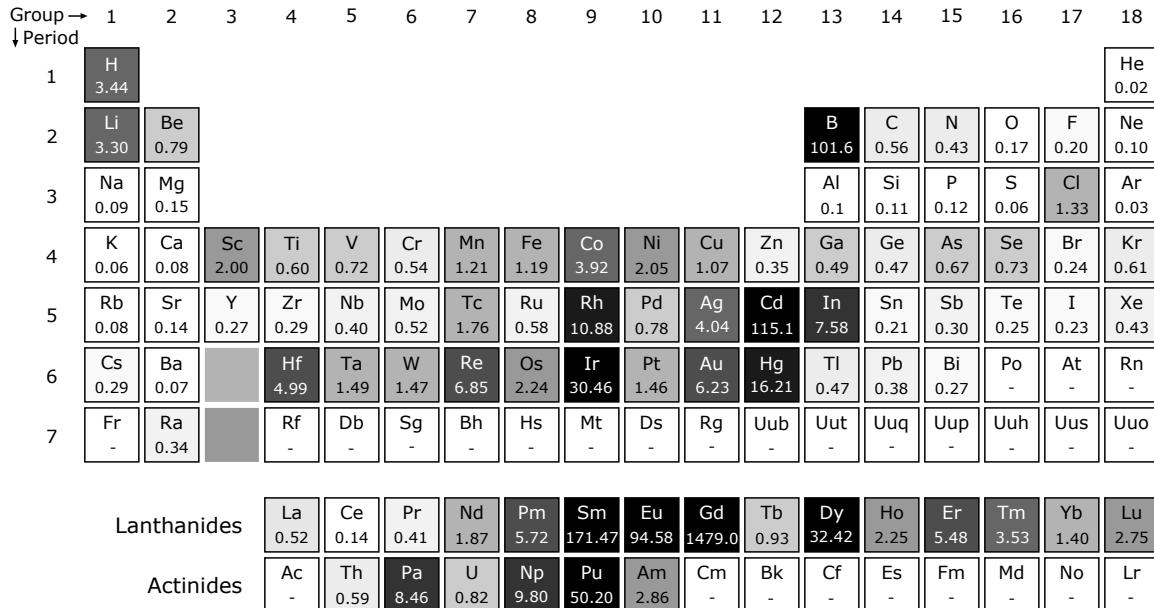


Fig. 2.5: Attenuation coefficients for thermal neutrons.

| Group → ↓ Period | 1 | 2 | 3 | 4 | 5 | 6 | 7 | 8 | 9 | 10 | 11 | 12 | 13 | 14 | 15 | 16 | 17 | 18 |
|------------------|------------|-------------|-------------|-------------|-------------|------------|-------------|-------------|-------------|-------------|-------------|-------------|-------------|-------------|-------------|-------------|------------|------------|
| 1 | H 0.02 | | | | | | | | | | | | | | | | He 0.02 | |
| 2 | Li 0.06 | Be 0.22 | | | | | | | | | | | | | | B 0.28 | C 0.27 | N 0.11 |
| 3 | Na 0.13 | Mg 0.24 | | | | | | | | | | | | | | Al 0.38 | Si 0.33 | P 0.25 |
| 4 | K 0.14 | Ca 0.26 | Sc 0.48 | Ti 0.73 | V 1.04 | Cr 1.29 | Mn 1.32 | Fe 1.57 | Co 1.78 | Ni 1.96 | Cu 1.97 | Zn 1.64 | Ga 1.42 | Ge 1.33 | As 1.50 | Se 1.23 | Br 0.90 | Kr 0.73 |
| 5 | Rb 0.47 | Sr 0.86 | Y 1.61 | Zr 2.47 | Nb 3.43 | Mo 4.29 | Tc 5.06 | Ru 5.71 | Rh 6.08 | Pd 6.13 | Ag 5.67 | Cd 4.84 | In 4.31 | Sn 3.98 | Sb 4.28 | Te 4.06 | I 3.45 | Xe 2.53 |
| 6 | Cs 1.47 | Ba 2.73 | | Hf 19.70 | Ta 25.47 | W 30.49 | Re 34.47 | Os 37.92 | Ir 39.01 | Pt 38.61 | Au 35.94 | Hg 25.88 | Tl 23.23 | Pb 22.81 | Bi 20.28 | Po 20.22 | At - | Rn 9.77 |
| 7 | Fr - | Ra 11.80 | | Rf - | Db - | Sg - | Bh - | Hs - | Mt - | Ds - | Rg - | Uub - | Uut - | Uuq - | Uup - | Uuh - | Uus - | Uuo - |
| Lanthanides | | La 5.04 | Ce 5.79 | Pr 6.23 | Nd 6.46 | Pm 7.33 | Sm 7.68 | Eu 5.66 | Gd 8.69 | Tb 9.46 | Dy 10.17 | Ho 10.17 | Er 11.70 | Tm 12.49 | Yb 9.32 | Lu 14.07 | | |
| Actinides | | Ac 24.47 | Th 28.95 | Pa 39.65 | U 49.08 | Np - | Pu - | Am - | Cm - | Bk - | Cf - | Es - | Fm - | Md - | No - | Lr - | | |

Fig. 2.6: Attenuation coefficients for X-rays.

2.5 Measurements are rarely perfect

There is no perfect measurement. This is also true for neutron imaging. The ideal image is the sample is distorted for many reasons. The figure below shows how an image of a sample can look after passing through the acquisition system. These quality degradations will have an impact on the analysis of the image data. In some cases, it is possible to correct for some of these artifacts using classic image processing techniques. There are however also cases that require extra effort to correct the artifacts.

2.5.1 Factors affecting the image quality

The list below provides some factors that affect the quality of the acquired images. Most of them can be handled by changing the imaging configuration in some sense. It can however be that the sample or process observed put limitations on how much the acquisition can be tuned to obtain the perfect image.

- Resolution (Imaging system transfer functions)
- Noise
- Contrast
- Inhomogeneous contrast
- Artifacts

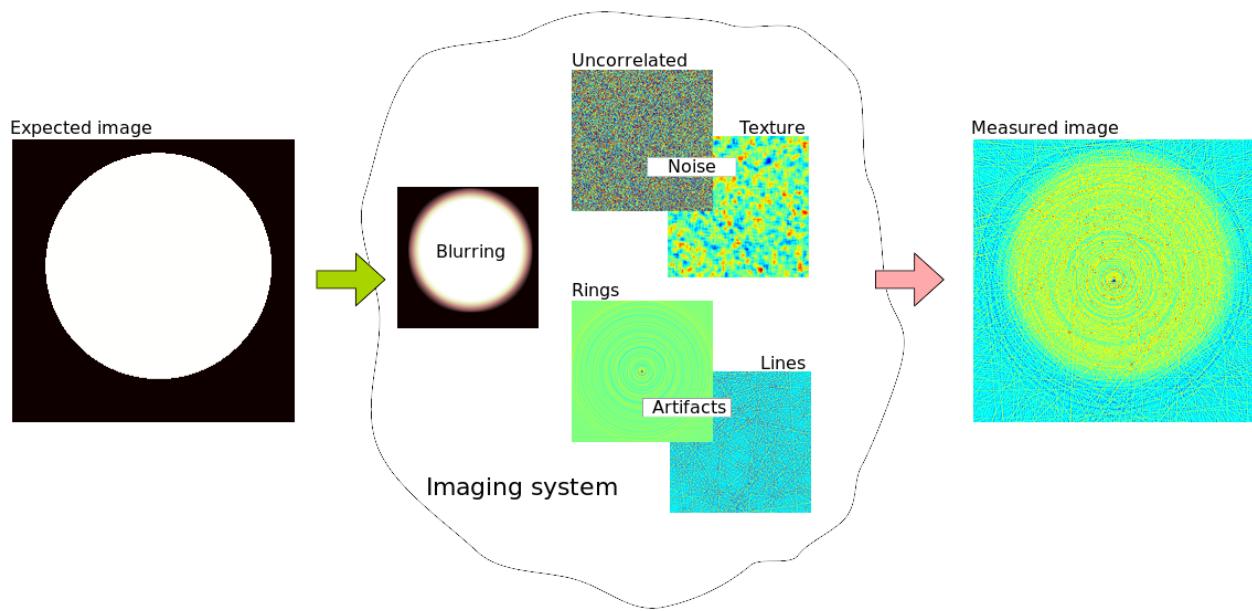


Fig. 2.7: Schematic showing the error contributions in an imperfect imaging system.

Resolution

The resolution is primarily determined by the optical transfer function of the imaging system. The actual resolution is given by the extent of the sample and how much the detector needs to capture in one image. This gives the field of view and given the number of pixels in the used detector it is possible to calculate the pixel size. The pixel size limits the size of the smallest feature in the image that can be detected. The scintillator, which is used to convert neutrons into visible light, is chosen to

1. match the sampling rate given by the pixel size.
2. provide sufficient neutron capture to obtain sufficient light output for a given exposure time.

Noise

An imaging system has many noise sources, each with its own distribution e.g.

1. Neutron statistics - how many neutrons are collected in a pixel. This noise is Poisson distributed.
2. Photon statistics - how many photons are produced by each neutron. This noise is also Poisson distributed.
3. Thermal noise from the electronics which has a Gaussian distribution.
4. Digitization noise from converting the charges collected for the photons into digital numbers that can be transferred and stored by a computer, this noise has a binomial distribution.

The neutron statistics are mostly dominant in neutron imaging but in some cases it could also be that the photon statistics play a role.

Contrast

The contrast in the sample is a consequence of

1. how well the sample transmits the chosen radiation type. For neutrons you obtain good contrast from materials containing hydrogen or lithium while many metals are more transparent.
2. the amount of a specific element or material represented in a unit cell, e.g. a pixel (radiograph) or a voxel (tomography).

The objective of many experiments is to quantify the amount of a specific material. This could for example be the amount of water in a porous medium.

Good contrast between different image features is important if you want to segment them to make conclusions about the image content. Therefore, the radiation type should be chosen to provide the best contrast between the features.

Inhomogeneous contrast

The contrast in the raw radiograph depends much on the beam profile. These variations are easily corrected by normalizing the images by an open beam or flat field image.

- **Biases introduced by scattering** Scattering is the dominant interaction for many materials used in neutron imaging. This means that neutrons that are not passing straight through the sample are scattered and contribute to a background cloud of neutrons that build up a bias of neutrons that are also detected and contribute to the image.
- **Biases from beam hardening** is a problem that is more present in x-ray imaging and is caused by the fact that the attenuation coefficient depends on the energy of the radiation. Higher energies have lower attenuation coefficient, thus will high energies penetrate the thicker samples than lower energies. This can be seen when a polychromatic beam is used.

Artifacts

Many images suffer from outliers caused by stray rays hitting the detector. Typical artefacts in tomography data are

- Lines, which are caused by outlier spots that only appear in single projections. These spots appear as lines in the reconstructed images.
- Rings are caused by stuck pixels which have the same value in all projections.

2.6 Introduction to segmentation

The basic task of segmentation is to identify regions of pixels with similar properties. This can be on different levels of abstraction. The lowest level is to create segments of pixels based on their pixel values like in the LANDSAT image below.

This type of segmentation sometimes be done with the help of a histogram that shows the distribution of values in the image.

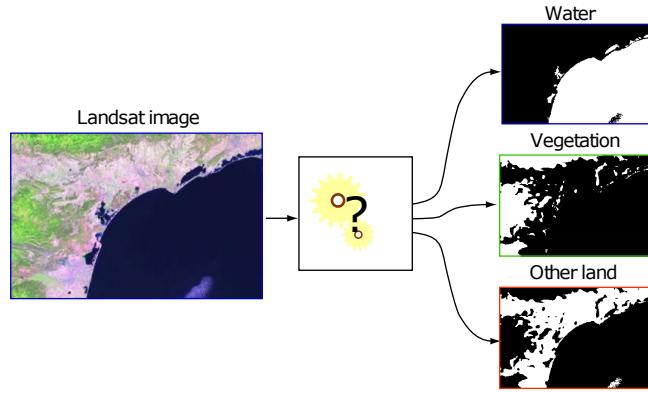
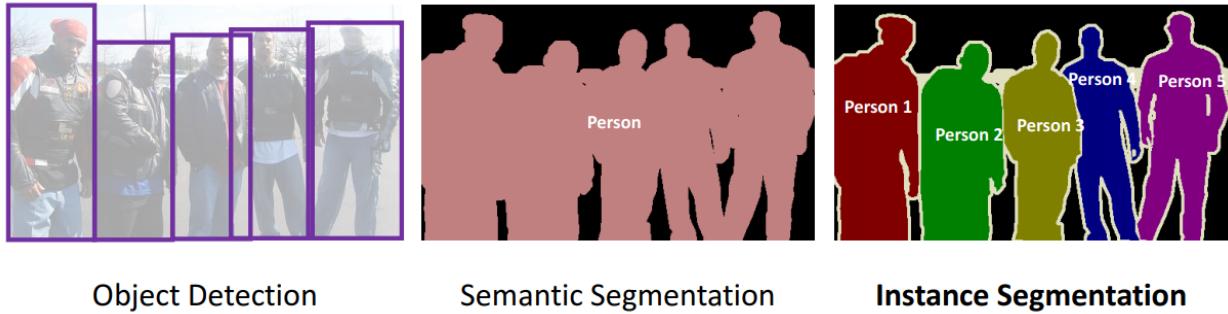


Fig. 2.8: Segmentation of a satellite image to identify different land types.

2.7 Different types of segmentation

When we talk about image segmentation there are different meanings to the word. In general, segmentation is an operation that marks up the image based on pixels or pixel regions. This is a task that has been performed since beginning of image processing as it is a natural step in the workflow to analyze images - we must know which regions we want to analyze and this is a tedious and error prone task to perform manually. Looking at the figure below we two type of segmentation.



Object Detection

Semantic Segmentation

Instance Segmentation

- Object detection - identifies regions containing an object. The exact boundary does not matter so much here. We are only interested in a bounding box.
- Semantic segmentation - classifies all the pixels of an image into meaningful classes of objects. These classes are “semantically interpretable” and correspond to real-world categories. For instance, you could isolate all the pixels associated with a cat and color them green. This is also known as dense prediction because it predicts the meaning of each pixel.
- Instance segmentation - Identifies each instance of each object in an image. It differs from semantic segmentation in that it doesn’t categorize every pixel. If there are three cars in an image, semantic segmentation classifies all the cars as one instance, while instance segmentation identifies each individual car.

2.7.1 Basic segmentation: Applying a threshold to an image

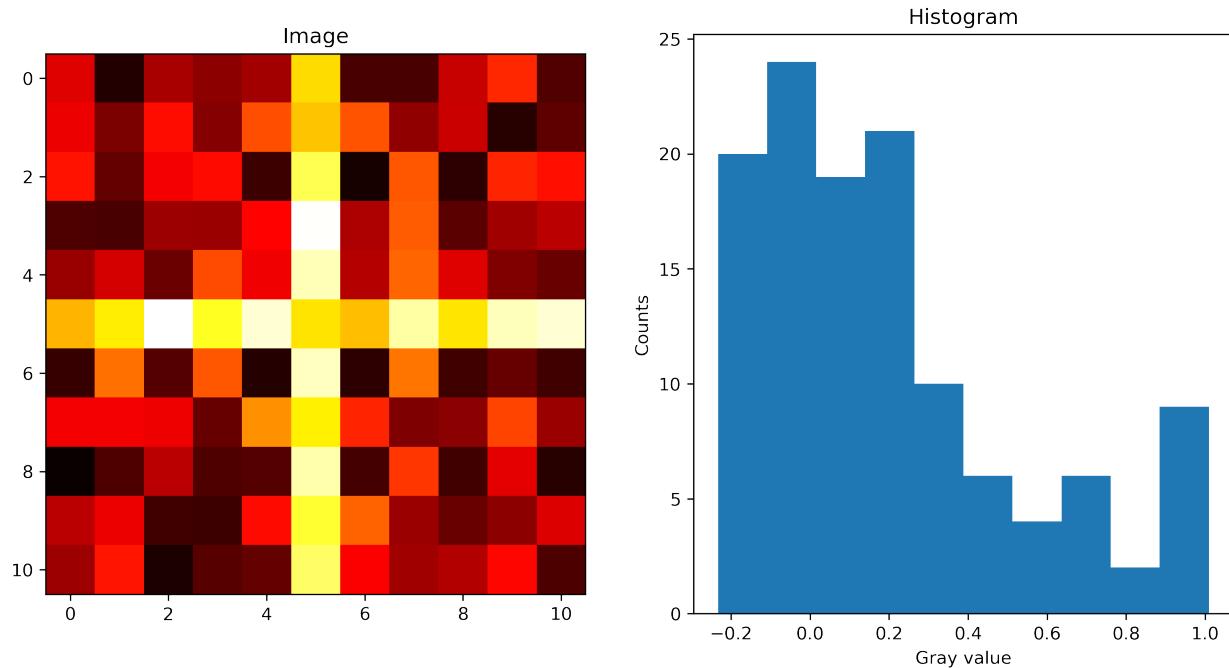
Start out with a simple image of a cross with added noise

$$I(x, y) = f(x, y)$$

Here, we create a test image with two features embedded in uniform noise; a cross with values in the order of ‘1’ and background with values in the order ‘0’. The figure below shows the image and its histogram. The histogram helps us to see how the graylevels are distributed which guides the decision where to put a threshold that segments the cross from the background.

```
fig,ax = plt.subplots(1,2,figsize=(12,6))
nx = 5; ny = 5;
# Create the test image
xx, yy = np.meshgrid(np.arange(-nx, nx+1)/nx*2*np.pi, np.arange(-ny, ny+1)/ny*2*np.pi)
cross_im = 1.5*np.abs(np.cos(xx*yy)) / (np.abs(xx*yy)+(3*np.pi/nx)) + np.random.uniform(-0.25, 0.25, size = xx.shape)

# Show it
im=ax[0].imshow(cross_im, cmap = 'hot'); ax[0].set_title("Image")
ax[1].hist(cross_im.ravel(),bins=10); ax[1].set_xlabel('Gray value'); ax[1].set_ylabel('Counts'); ax[1].set_title("Histogram");
```



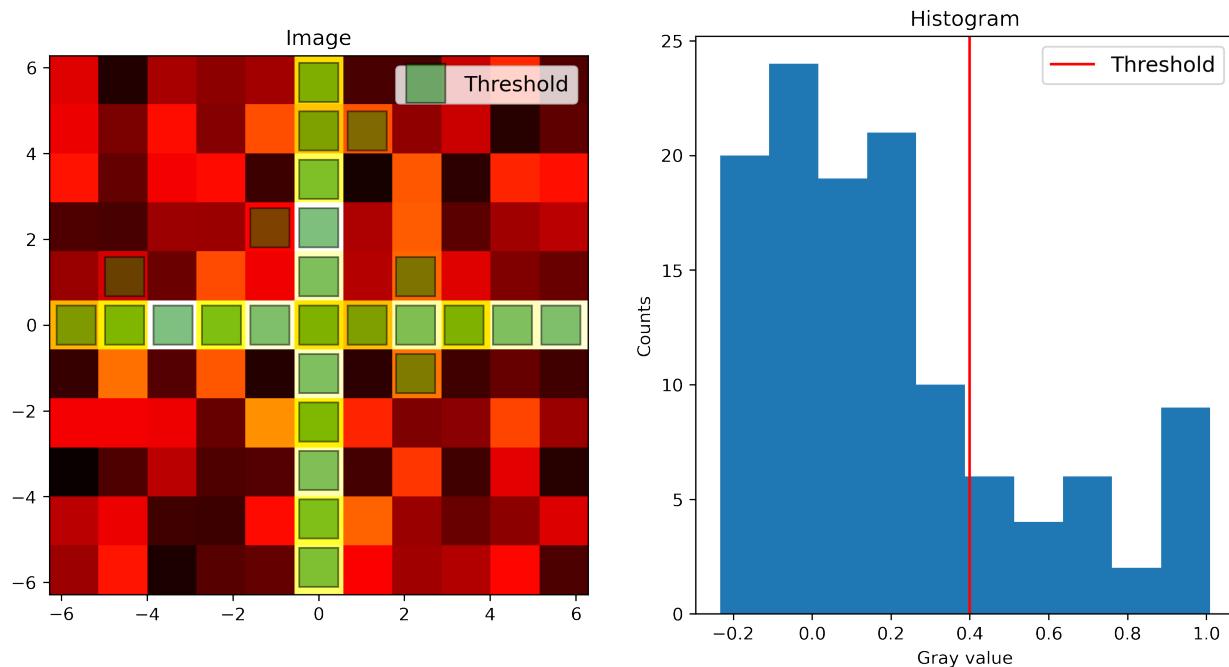
2.7.2 Applying a threshold to an image

By examining the image and probability distribution function, we can *deduce* that the underlying model is a whitish phase that makes up the cross and the darkish background

Applying the threshold is a deceptively simple operation

$$I(x, y) = \begin{cases} 1, & f(x, y) \geq 0.40 \\ 0, & f(x, y) < 0.40 \end{cases}$$

```
threshold = 0.4; thresh_img = cross_im > threshold
fig,ax = plt.subplots(1,2,figsize=(12, 6))
ax[0].imshow(cross_im, cmap = 'hot', extent = [xx.min(), xx.max(), yy.min(), yy.
    max()]); ax[0].set_title("Image")
ax[0].plot(xx[np.where(thresh_img)]*0.9, yy[np.where(thresh_img)]*0.9,
    'ks', markerfacecolor = 'green', alpha = 0.5,label = 'Threshold',_
    markeredgecolor='black', markersize = 22); ax[0].legend(fontsize=12);
ax[1].hist(cross_im.ravel(),bins=10); ax[1].axvline(x=threshold,color='r',label=
    'Threshold'); ax[1].legend(fontsize=12);
ax[1].set_xlabel('Gray value'); ax[1].set_ylabel('Counts'); ax[1].set_title("Histogram
    "));
```



In this fabricated example we saw that thresholding can be a very simple and quick solution to the segmentation problem. Unfortunately, real data is often less obvious. The features we want to identify for our quantitative analysis are often obscured by different other features in the image. They may be part of the setup or caused by the acquisition conditions.

2.8 Noise and SNR

Any measurement has a noise component and this noise has to be dealt with in some way when the image is to be segmented. When the noise is not handled correctly it will cause many misclassified pixels. The noise strength of measured using the signal to noise ratio **SNR**. It is defined as the ratio between signal average and signal standard deviation.

The noise in neutron imaging mainly originates from the amount of captured neutrons.

This noise is Poisson distributed and the signal to noise ratio is

$$SNR = \frac{E[x]}{s[x]} \sim \frac{N}{\sqrt{N}} = \sqrt{N}$$

where N is the number of captured neutrons. The figure below shows two neutron images acquired at 0.1s and 10s respectively. The plot shows the signal to noise ratio obtained for different exposure times.

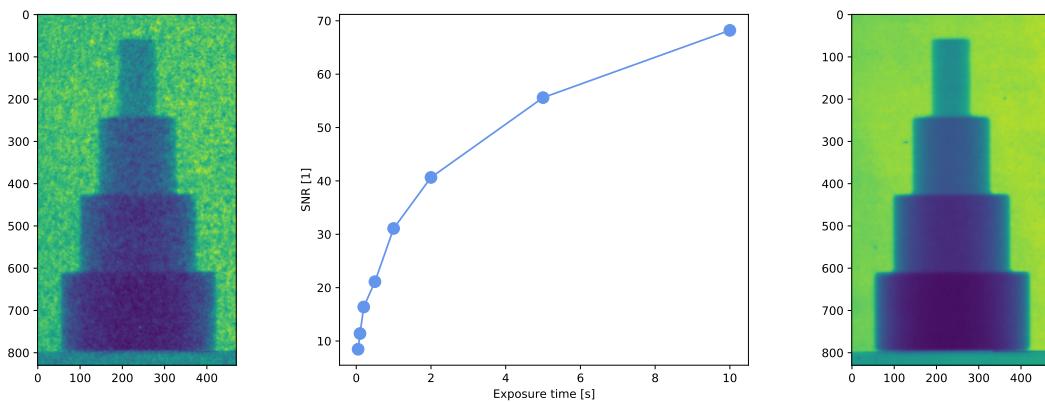


Fig. 2.9: Signal to noise ratio for radiographies acquired with different exposure times.

The signal to noise ratio can be improved by increasing the number of neutrons per pixel. This can be achieved through increasing

- Neutron flux - this is usually relatively hard as the neutron sources operate with the parameters it is designed for. There is a possibility by changing the neutron aperture, but has an impact of the beam quality.
- Exposure time - the exposure time can be increased but in the end there is a limitation on how much this can be used. Beam time is limited which means the experiment must be finished in a given time. There is also an upper limit on the exposure time defined by the observed sample or process when it changes over time. Too long exposure times will result in motion artefacts.
- Pixel size - increasing the pixel size means that neutrons are collected over a greater area and thus more neutrons are captured during the exposure. The limit on how much you can increase the pixel size is defined by the smallest features you want to detect.
- Detector material and thickness - the number of captured neutrons depends on the scintillator material and how thick it is. The thickness does however have an impact on the resolution. Therefore scintillator thickness and pixel size often increase in parallel as there is no point in oversampling a smooth signal to much.

In the end, there are many parameters that combined results in the SNR you obtain. These parameters are tuned to match the experiment conditions. Filtering techniques can help to increase the SNR still it is too low for your quantitative analysis.

2.9 Problematic segmentation tasks

Segmentation is rarely an obvious task. What you want to find is often obscured by other image features and unwanted artefacts from the experiment technique. If you take a glance at the painting by Bev Doolittle, you quickly spot the red fox in the middle. Looking a little closer at the painting, you'll recognize two Indians on their spotted ponies. This example illustrates the problems you will encounter when you start to work with image segmentation.



Fig. 2.10: Cases making the segmentation task harder than just applying a single threshold.

Woodland Encounter Bev Doolittle

2.9.1 Typical image features that makes life harder

The basic segmentation shown in the example above can only be used under good conditions when the classes are well separated. Images from many experiments are unfortunately not well-behaved in many cases. The figure below shows four different cases when an advanced technique is needed to segment the image. Neutron images often show a combination of all cases.

The impact on the segmentation performance of all these cases can be reduced by proper experiment planning. Still, sometimes these improvements are not fully implemented in the experiment to be able to fulfill other experiment criteria.

In neutron imaging you see all these image phenomena.

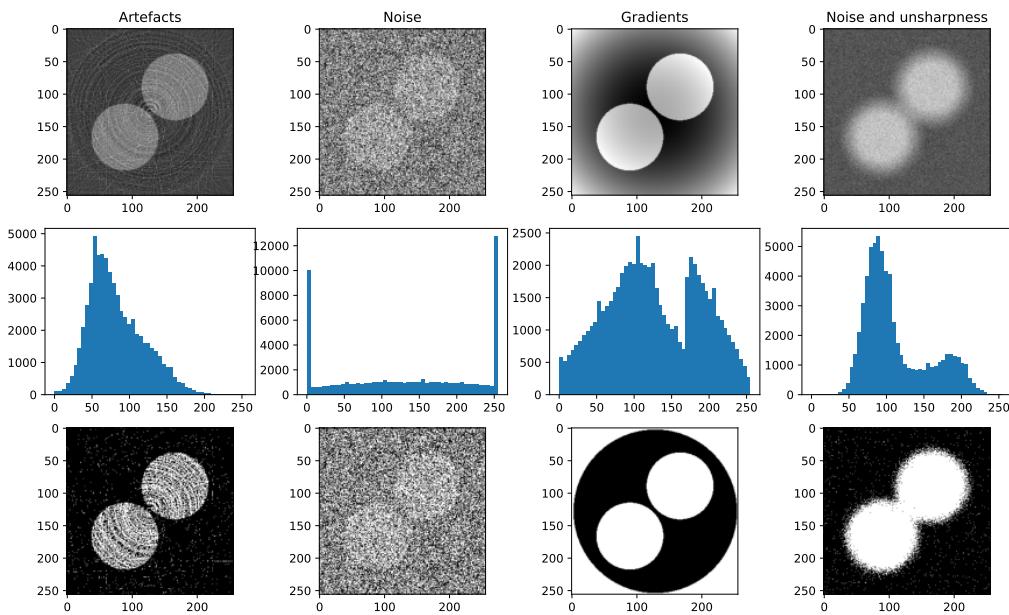


Fig. 2.11: Cases making the segmentation task harder than just applying a single threshold.

2.10 Segmentation problems in neutron imaging

LIMITED DATA PROBLEM

Machine learning methods require a lot of training data to be able to build good models that are able to detect the features they are intended to.

Different types of limited data:

- Few data points or limited amounts of images

This is very often the case in neutron imaging. The number of images collected during an experiment session is often very low due to the long experiment duration and limited amount of beam time. This makes it hard to develop segmentation and analysis methods for single experiments. The few data points problem can partly be overcome by using data from previous experiment with similar characteristics. The ability to recycle data depends on what you want to detect in the images.

- Unbalanced data

Unbalanced data means that the ratio between the data points with features you want detect and the total number data points is several orders of magnitude. E.g roots in a volume like the example we will look at later in this lecture. There is even a risk that the distribution of the wanted features is overlapped by the dominating background distribution.

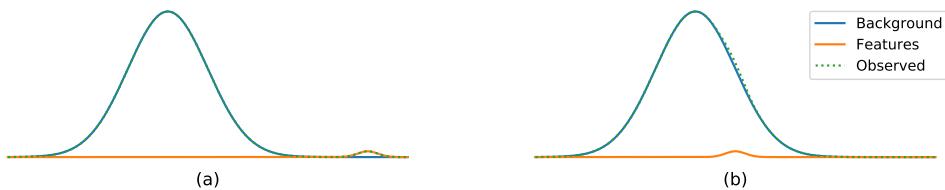


Fig. 3.1: Two cases of unbalanced data; (a) the classes are well separated and the feature class is clearly visible in the tail distribution of the background and (b) the feature class is embedded in the background making it hard to detect.

Case (a) can most likely be segmented using one of the many histogram based thresholding methods proposed in literature. Case (b) is much harder to segment as the target features have similar gray levels as the background. This case requires additional information to make segmentation possible.

- Little or missing training data

A complete set of training data contains both input data and labelled data. The input data is easy to obtain, it is the images you measured during your experiment. The labelled data is harder to get as it is a kind of chicken and egg problem. In particular, if your experiment data is limited. In that case, you would have to mark-up most of the available data to obtain the labeled data. Which doesn't make sense because

- then you'd already solved the task before even starting to train your segmentation algorithm.
- An algorithm based on learning doesn't improve the results, it only make it easier to handle large amounts of data.

3.1 Training data from NI is limited

The introducing words about limited data essentially describes the problems that arise in neutron imaging. There are several reasons for this. Some are:

- Long experiment times
- Few samples
- Some recycling from previous experiments is possible.

The experiment times are usually rather long which results in only few data sets per experiment. In some cases there is the advantage that there are plenty experiments over the years that produced data with similar characteristics. These experiments can be used in a pool of data.

3.2 Augmentation to increase training data

Obtaining more experiment data is mostly relatively hard in neutron imaging. The available time is very limited. Still, many supervised analysis methods require large data sets to perform reliably. A method to improve this situation is to use data augmentation. The figure below shows some examples of augmentations of the same image. You can also add noise and modulate the image intensity to increase the variations further.

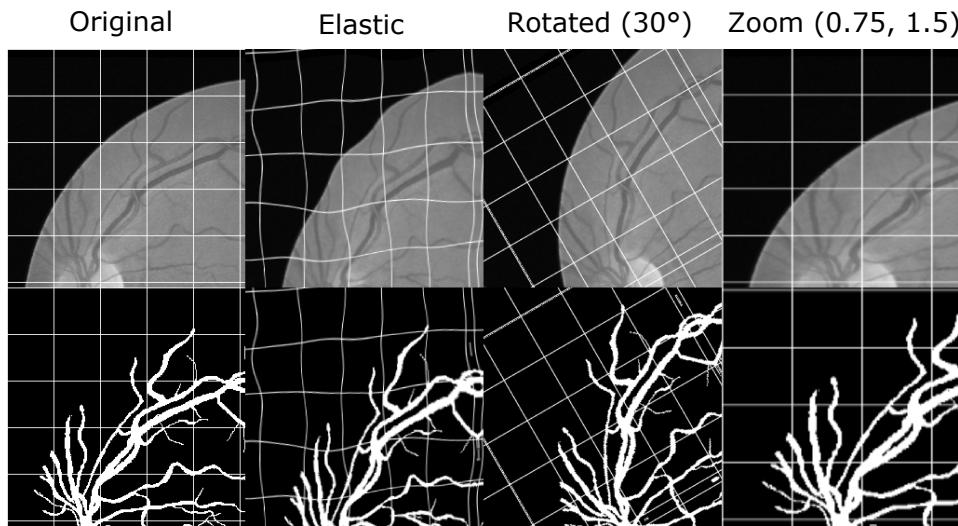


Fig. 3.2: A retinal image modified using different augmentation techniques (source: <https://drive.grand-challenge.org/DRIVE/>) prepared by Gian Guido Parenza.

Data augmentation is a method to modify your existing data to obtain variations of it.

Augmentation will be used to increase the training data in the root segmentation example in the end of this lecture.

3.3 Simulation to increase training data

A further way to increase training data is to build a model of the features you want to train on. This approach has the advantage that you know where to look for the features which means the tedious annotation task is reduced to a minimum. The work rather lies in building a reliable model that should reflect the characteristics of features you want to segments. Once a valid model is built, it is easy to generate masses of data under various conditions.

- Geometric models
- Template models
- Physical models

Both augmented and simulated data should be combined with real data.

3.4 Transfer learning

Many machine learning techniques also have the ability to remember what they learned before. This effect can be used to improve the performance also when little data is available from the current experiment.

Transfer learning is a technique that uses a pre-trained network to

- Speed up training on your current data
- Support in cases of limited data
- Improve network performance

Usually, the input and downsampling part of the network is used from the pre-trained network. It can be set as constant or trainable. The upsampling and output will be trained with new data from your current data set. Transfer learning will be covered in the root segmentation example in the end of this lecture.

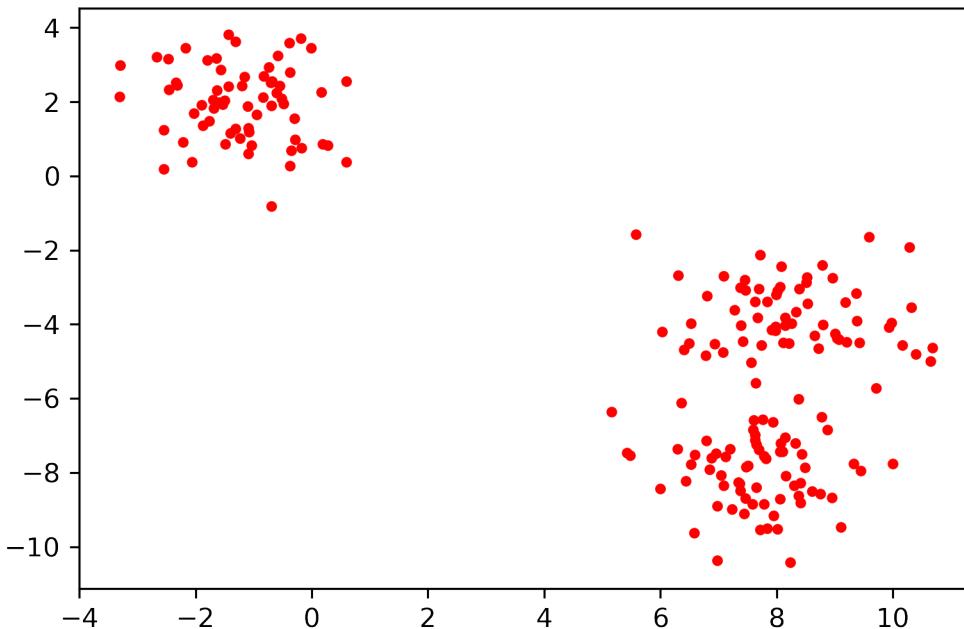
UNSUPERVISED SEGMENTATION

Unsupervised segmentation method tries to make sense of the that without any prior knowledge. They may need an initialization parameter telling how many classes are expected in the data, but beyond that you don't have to provide much more information.

4.1 Introducing clustering

With clustering methods you aim to group data points together into a limited number of clusters. Here, we start to look at an example where each data point has two values. The test data is generated using the `make_blobs` function.

```
test_pts = pd.DataFrame(make_blobs(n_samples=200, random_state=2018) [  
    0], columns=['x', 'y'])  
plt.plot(test_pts.x, test_pts.y, 'r.');
```



The generated data set has two obvious clusters, but if look closer it is even possible to identify three clusters. We will now use this data to try the k-means algorithm.

4.2 k-means

The k-means algorithm is one of the most used unsupervised clustering method.

It is an iterative method that starts with a label image where each pixel has a random label assignment. Each iteration involves the following steps:

1. Compute current centroids based on the o current labels
2. Compute the value distance for each pixel to the each centroid. Select the class which is closest.
3. Reassign the class value in the label image.
4. Repeat until no pixels are updated.

The distance from pixel i to centroid j is usually computed as $\|p_i - c_j\|_2$.

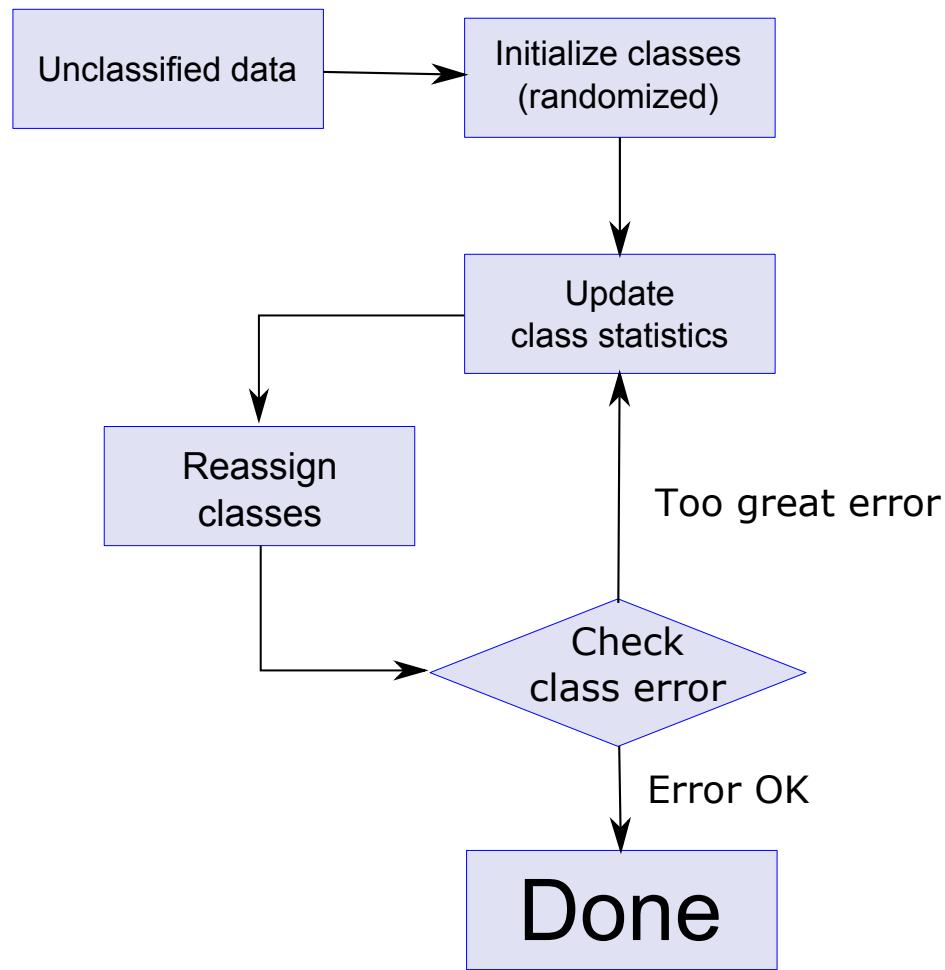


Fig. 4.1: Flow chart for the k-means clustering iterations.

The user only have to provide the number of classes the algorithm shall find.

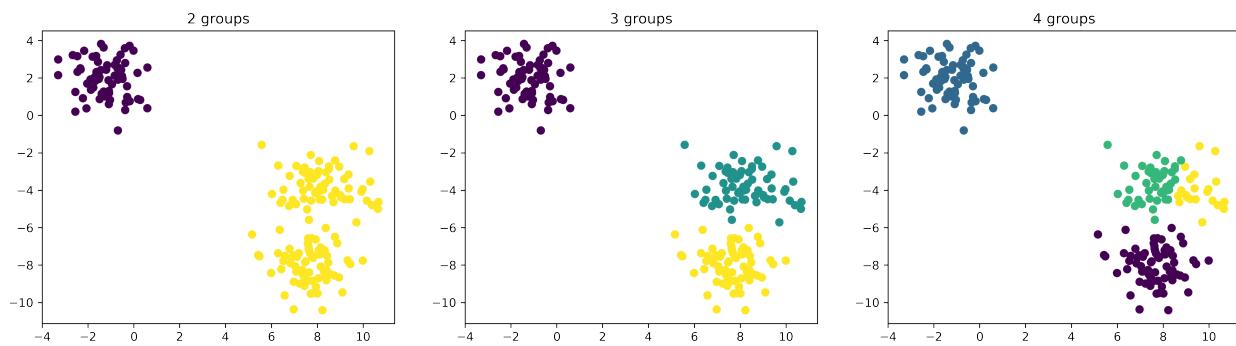
Note The algorithm will find exactly the number you ask it to, it doesn't care if it makes sense!

4.3 Basic clustering example

In this example we will use the blob data we previously generated and to see how k-means behave when we select different numbers of clusters.

```
N=3
fig, ax = plt.subplots(1,N,figsize=(18, 4.5))

for i in range(N) :
    km = KMeans(n_clusters=i+2, random_state=2018); n_grp = km.fit_predict(test_pts)
    ax[i].scatter(test_pts.x, test_pts.y, c=n_grp)
    ax[i].set_title('{0} groups'.format(i+2))
```

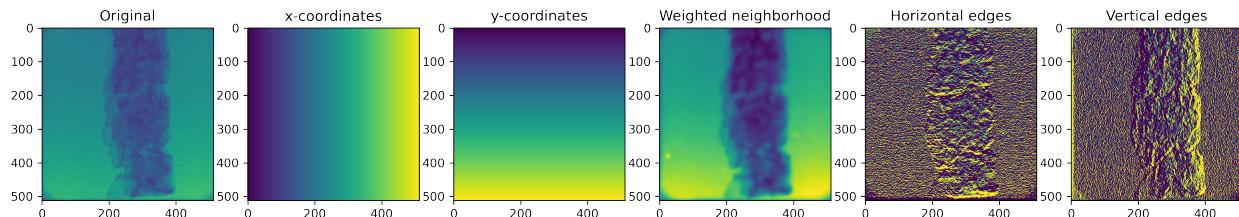


When we select two clusters there is a natural separation between the two clusters we easily spotted by just looking at the data. When the number of clusters is increased to three, we again see a cluster separation that makes sense. Now, when the number of clusters is increased yet another time we see that one of the clusters is split once more. This time it is however questionable if the number of clusters makes sense. From this example, we see that it is important to be aware of problems related to over segmentation.

4.4 Add spatial information to k-means

It is important to note that k-means by definition is not position sensitive. Clustering is by definition not position sensitive, mainly measures distances between values and distributions. The position can however be included as additional components in the data vectors. You can also add neighborhood information using filtered images as additional components of the feature vectors.

```
orig = fits.getdata('../data/spots/mixture12_00001.fits')[:, :, ::4]
fig, ax = plt.subplots(1, 6, figsize=(18, 5)); x, y = np.meshgrid(np.linspace(0, 1, orig.shape[0]), np.linspace(0, 1, orig.shape[1]))
ax[0].imshow(orig, vmin=0, vmax=4000), ax[0].set_title('Original')
ax[1].imshow(x), ax[1].set_title('x-coordinates')
ax[2].imshow(y), ax[2].set_title('y-coordinates')
ax[3].imshow(flt.gaussian(orig, sigma=5)), ax[3].set_title('Weighted neighborhood')
ax[4].imshow(flt.sobel_h(orig), vmin=0, vmax=0.001), ax[4].set_title('Horizontal edges')
ax[5].imshow(flt.sobel_v(orig), vmin=0, vmax=0.001), ax[5].set_title('Vertical edges');
```



Scaling the values is very important when you use these additional feature vectors. Otherwise, some vectors will dominate others.

4.5 When can clustering be used on images?

Clustering and in particular k-means are often used for imaging applications.

- Single images

It does not make much sense to segment single images using k-means. This would result in thresholding similar to the one provided by Otsu's method. If you, however, add spatial information like edges and positions it starts become interesting to use k-means for a single image.

- Bimodal data

In recent years, several neutron imaging instruments have installed an X-ray source to provide complementary information to the neutron images. This is a great mix of information for k-means based segmentation. Another type of bimodal data is when a grating interferometry setup is used. This setup provides three images revealing different aspects of the samples. Again, here it is a great combination as these data are

- Spectrum data

Many materials have a characteristic response to the neutron wavelength. This is used in many materials science experiments, in particular experiments performed at pulsed neutron sources. The data from such experiments result in a spectrum response for each pixel. This means each pixel can be a vector of >1000 elements.

4.6 Clustering applied to wavelength resolved imaging

4.6.1 The imaging techniques and its applications

4.6.2 The data

In this example we use a time of flight transmission spectrum from an assembly of six rectangular steel blocks produced by means of additive manufacturing. The typical approach to analyse this kind of information is to select some regions of interest and compute the average spectra from these regions. The average spectra are then analyzed using single Bragg edge fitting methods or Rietveld refinement.

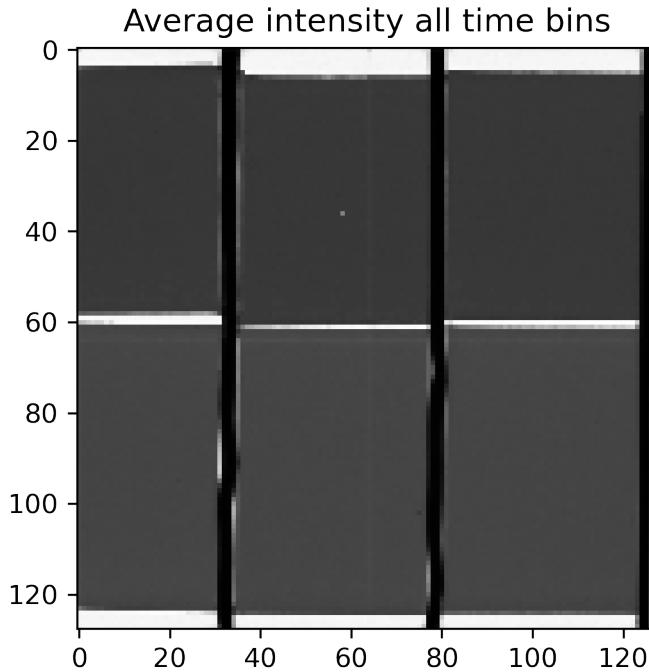
Here, we will explore the possibility to identify regions with similar spectra using k-means clustering. The data is provided on the repository and has the dimensions 128x128x661, where the first two dimensions reflect the image size and the last is the number of time bins of the time of flight spectrum. This is a downsized version of the original data which has 512x512 image frames. Each frame has a relatively low signal to noise ratio, therefore we use the average image `wtof` to show the image contents.

```
toto = np.load('../data/todata.npy')
wtof = toto.mean(axis=2)
```

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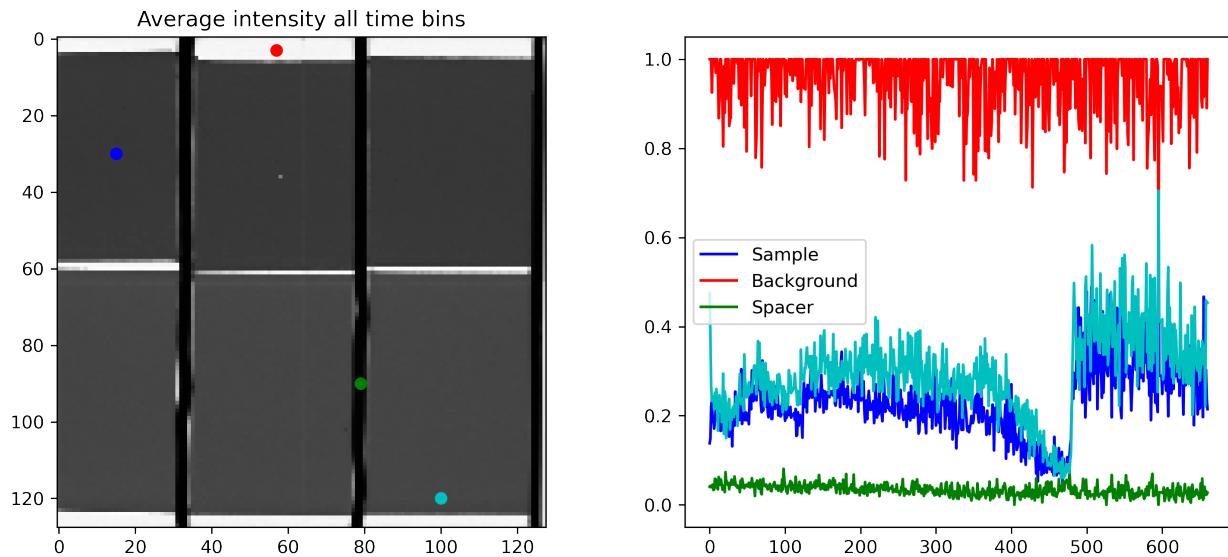
```
plt.imshow(wtof, cmap='gray');
plt.title('Average intensity all time bins');
```



4.6.3 Looking at the spectra

The average image told us that there are six sample regions, some void, and some spacers separating the samples. Now you might think, “easy” just apply some thresholds on the image and you have identified these regions. Now, the pixels in this image are represented by over 600 spectrum samples which reveal a different story. The sample was exposed to a surface treatment and we expect to locate regions within the samples. Given this information it is much harder to identify the regions of interest.

```
fig, ax= plt.subplots(1,2,figsize=(12,5))
ax[0].imshow(wtof,cmap='gray'); ax[0].set_title('Average intensity all time bins');
ax[0].plot(57,3,'ro'), ax[0].plot(15,30,'bo'), ax[0].plot(79,90,'go'); ax[0].plot(100,
˓→120,'co');
ax[1].plot(tof[30,15,:],'b', label='Sample'); ax[1].plot(tof[3,57,:],'r', label=
˓→'Background'); ax[1].plot(tof[90,79,:],'g', label='Spacer'); ax[1].legend(); ax[1].
˓→plot(tof[120,100,:],'c', label='Sample 2');
```



These plots are very noisy and it may even be hard to identify regions with the same spectrum. This is where k-means may come to you help.

Reshaping

The k-means requires vectors for each feature dimension, not images as we have in this data set. Therefore, we need to reshape our data. The reshaping is best done using the numpy array method `reshape` like this:

```
tofr=tot.reshape([tot.shape[0]*tot.shape[1],tot.shape[2]])
print("Input ToF dimensions",tot.shape)
print("Reshaped ToF data",tofr.shape)
```

```
Input ToF dimensions (128, 128, 661)
Reshaped ToF data (16384, 661)
```

The reshaped data `tofr` now has the dimensions 16384x661, i.e. each 128x128 pixel image has been replaced by a vector with the length 16384 elements. The number of time bins still remains the same. The `reshape` method is a cheap operation as it only changes the elements of the dimension vector and doesn't rearrange the data in memory.

4.6.4 Setting up and running k-means

When we set up k-means, we merely have to select the number of clusters. The choice depends on many factors. Particularly in this case with data points containing a large number of values we have a great degree of freedom in what could be considered the correct result. Thus we should ask ourselves what we expect from the clustering.

- We can clearly see that there is void on the sides of the specimens.
- There is also a separating band between the specimens.
- Finally we have to decide how many regions we want to find in the specimens. Let's start with two regions with different characteristics.

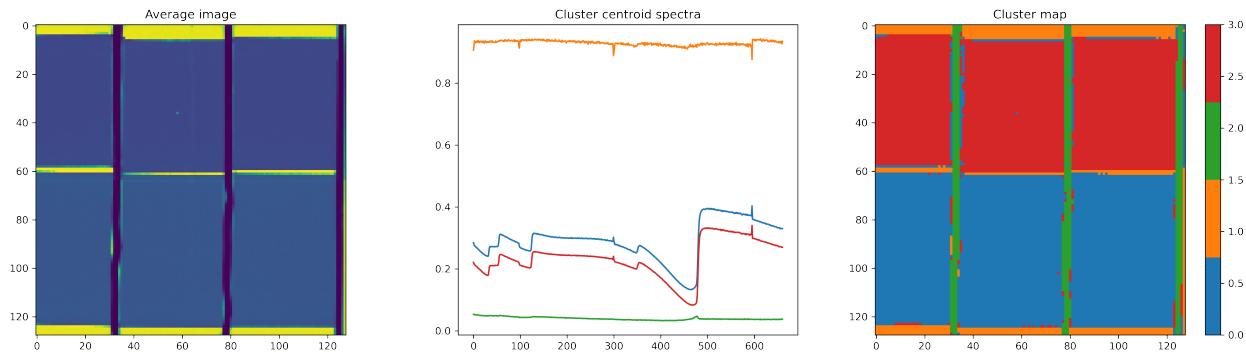
```
km = KMeans(n_clusters=4, random_state=2018)      # Random state is an initialization
# parameter for the random number generator
c = km.fit_predict(tofr).reshape(tof.shape[:2])   # Label image
kc = km.cluster_centers_.transpose()               # cluster centroid spectra
```

Results from the first try

```

fig,axes = plt.subplots(1,3,figsize=(18,5)); axes=axes.ravel()
axes[0].imshow(wtof,cmap='viridis'); axes[0].set_title('Average image')
p=axes[1].plot(kc);
axes[1].set_title('Cluster centroid spectra');
axes[1].set_aspect(tof.shape[2], adjustable='box')
cmap=ps.buildCMap(p) # Create a color map with the same colors as the plot

im=axes[2].imshow(c,cmap=cmap); plt.colorbar(im);
axes[2].set_title('Cluster map');
plt.tight_layout()
    
```



An advantage of the k-means is that it computes the centroid spectra for you as you go along. This means that if you are interested in finding average spectra in the identified regions, you don't need to compute them as a postprocessing step. They are already there!

Now, let's look at the spectra we get from the four-cluster segmentation. Two clusters are trivial; void and spacers. The remaining two clusters show us that there are two types of samples in this assembly... but they don't show if there is an effect of the surface treatments. These effects are not visible in the segmented image. So, this result was not very convincing for our investigation...

4.6.5 We need more clusters

- Experiment data has variations on places we didn't expect k-means to detect as clusters.

Except for only showing the main features in the image we got some misclassified pixels. This happens in particular near the sample edges.

- We need to increase the number of clusters!

Increasing the number of clusters will maybe give us more regions than we actually are asking for, but the advantage is that we also get the regions we want to see.

Increasing the number of clusters

What happens when we increase the number of clusters to ten?

```

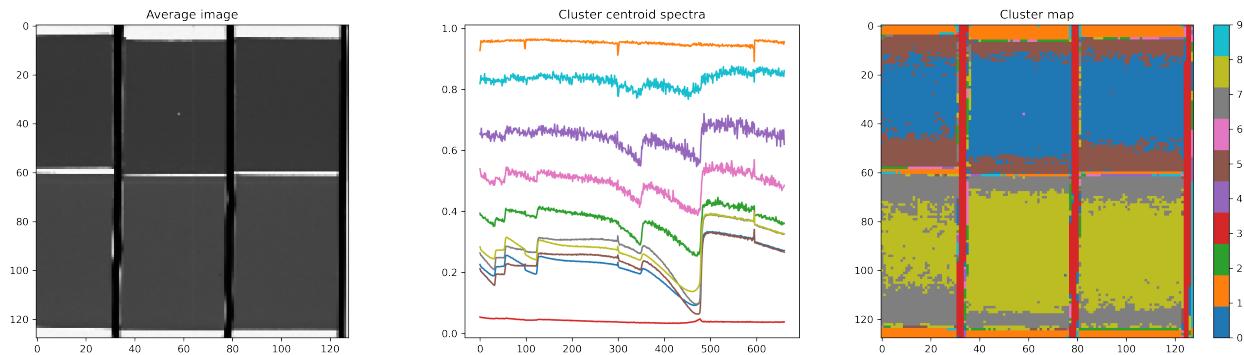
km = KMeans(n_clusters=10, random_state=2018)
c = km.fit_predict(tofr).reshape(tof.shape[:2]) # Label image
kc = km.cluster_centers_.transpose()           # cluster centroid spectra
    
```

Results of k-means with ten clusters

Segmenting neutron images using machine learning

```
fig,axes = plt.subplots(1,3,figsize=(18,5)); axes=axes.ravel()
axes[0].imshow(wtof,cmap='gray'); axes[0].set_title('Average image')
p=axes[1].plot(kc); axes[1].set_title('Cluster centroid spectra');
axes[1].set_aspect(tof.shape[2], adjustable='box')
cmap=ps.buildCMap(p) # Create a color map with the same colors as the plot

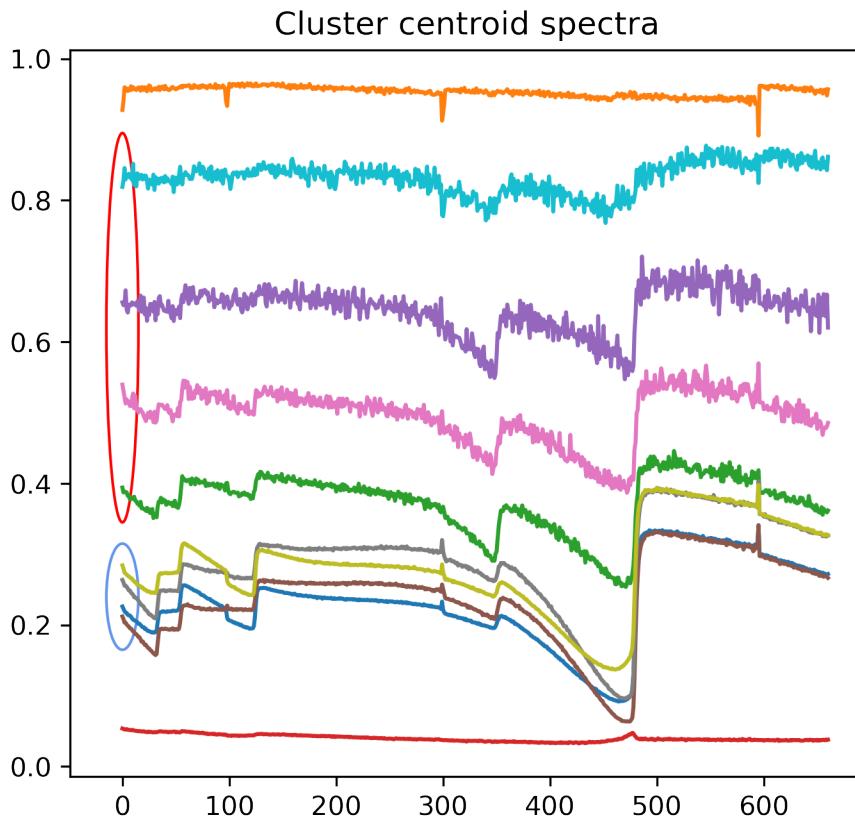
im=axes[2].imshow(c,cmap=cmap); plt.colorbar(im);
axes[2].set_title('Cluster map');
plt.tight_layout()
```



Interpreting the clusters

The segmentation with ten clusters did produce too many segments. If you look closer at these segments you see that they are mainly localized near the edges. This is a typical phenomenon in multi-class segmentation of images with smooth edges (transitions from one category to another). The red ellipse in the plot below highlights the centroids related to edge uncertainty.

```
fig,axes = plt.subplots(1,1,figsize=(14,5));
plt.plot(kc); axes.set_title('Cluster centroid spectra');
axes.add_patch(Ellipse((0,0.62), width=30,height=0.55,fill=False,color='r')) #,axes.
#set_aspect(tof.shape[2], adjustable='box')
axes.add_patch(Ellipse((0,0.24), width=30,height=0.15,fill=False,color='cornflowerblue
#)),axes.set_aspect(tof.shape[2], adjustable='box');
```



The blue ellipse shows the spectra which we are interested in. From these we can see the that there are actually two types of Bragg edge spectra represented in the six samples. These are transmission spectra and the variations in amplitude is related to sample thickness.

Cleaning up the works space

You may run short of memory after this part. You can delete the variables from this chapter to prepare yourself for the next chapter.

```
del km, c, kc, tofr, tof
```


SUPERVISED SEGMENTATION

The supervised method needs to be told how to behave when it is presented with some new data. This is done during a training session.

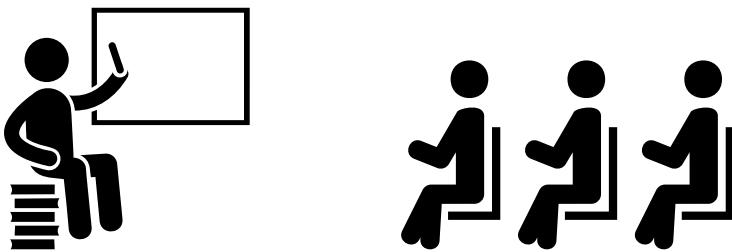


Fig. 5.1: Supervised methods require a training session before performing the actual segmentation task.

The typical workflow with a supervised method involves the following three steps:

1. **Training:** Requires training data
2. **Verification:** Requires verification data
3. **Inference:** The images you want to segment

The data for the training and verification steps need the same type of data, i.e., data with markup that tell which parts of the image belongs to which segment type. It is how even not good if you use the same data for both training and verification. That would be a self-fulfilling prophecy and the performance numbers for the trained model will be misleadingly high. Typically, you want to have more data for training than for verification. After all, the training data is the part that builds the model and you want good statistic foundation for the model.

5.1 k nearest neighbors

The *k nearest neighbors* algorithm makes the inference based on a point cloud of training point. When the model is presented with a new point it computes the distance to the closest points in the model. The *k* in the algorithm name indicates how many neighbors should be considered. E.g. $k=3$ means that the major class of the three nearest neighbors is assigned the tested point. Looking at the example below we would say that using three neighbors the

- Green hiker would claim he is in a spruce forest.
- Orange hiker would claim he is in a mixed forest.
- Blue hiker would claim he is in a birch forest.

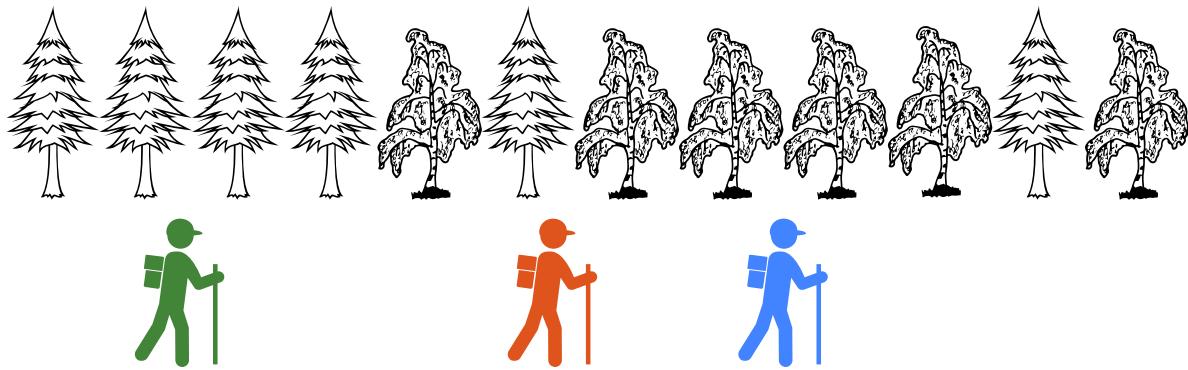
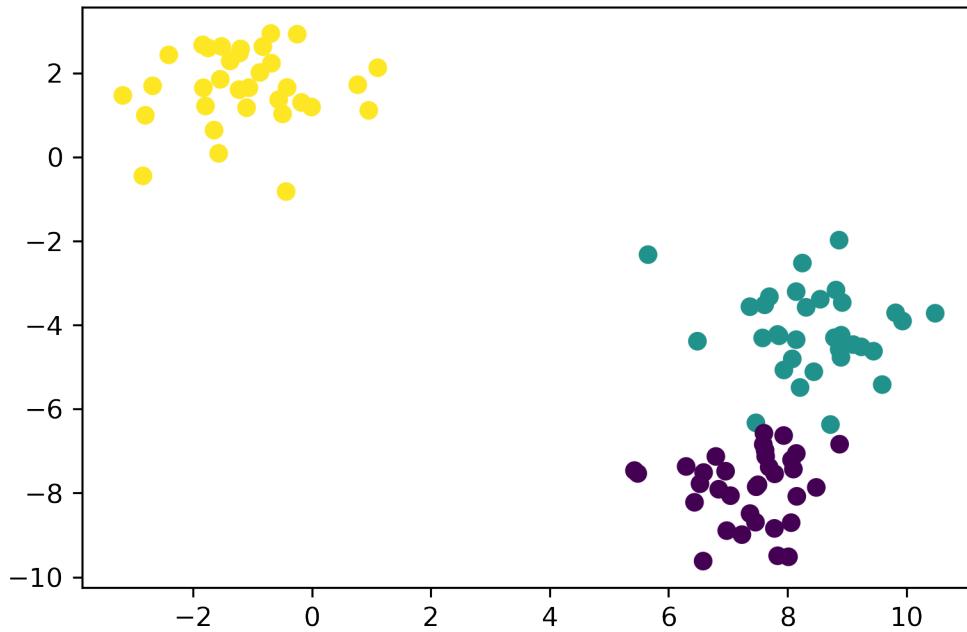


Fig. 5.2: Depending on the where the hiker is standing he makes the the conclusion that is either in a birch or spruce forest.

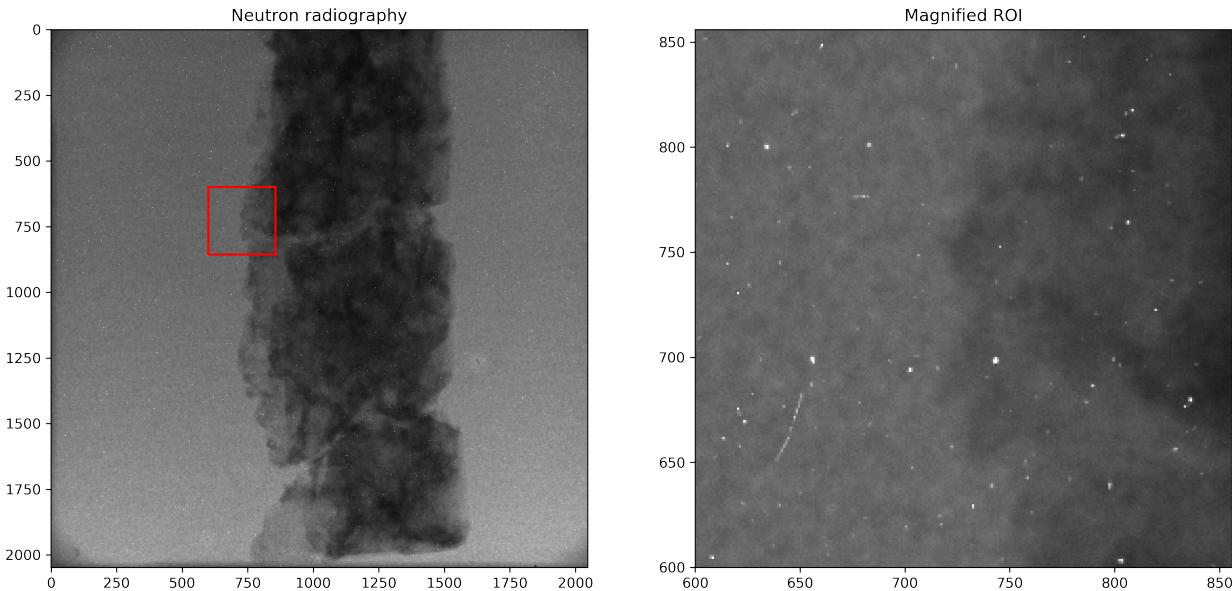
5.2 Create example data for supervised segmentation

```
blob_data, blob_labels = make_blobs(n_samples=100, random_state=2018)
test_pts = pd.DataFrame(blob_data, columns=['x', 'y'])
test_pts['group_id'] = blob_labels
plt.scatter(test_pts.x, test_pts.y, c=test_pts.group_id, cmap='viridis');
```



5.3 Detecting unwanted outliers in neutron images

```
orig= fits.getdata('..../data/spots/mixture12_00001.fits')
annotated=io.imread('..../data/spots/mixture12_00001.png'); mask=(annotated[:, :, 1]==0)
r=600; c=600; w=256
ps.magnifyRegion(orig,[r,c,r+w,c+w],[15,7],vmin=400,vmax=4000,title='Neutron
˓radiography')
```



5.3.1 Marked-up spots

5.3.2 Baseline - Traditional spot cleaning algorithm

Parameters

- N Width of median filter.
- k Threshold level for outlier detection.

5.3.3 Bivariate histogram of the detection image

```
selem=np.ones([3,3])
forig=orig.astype('float32')
mimg = flt.median(forig,selem=selem)
d = np.abs(forig-mimg)

fig,ax=plt.subplots(1,2,figsize=(12,5))
h,x,y,u=ax[0].hist2d(forig.ravel(),d.ravel(), bins=100);
ax[0].set_xlabel('Input image - $f$'),ax[0].set_ylabel('$|f - \text{med}_{3x3}(f)|$'),ax[0].
˓set_title('Bivariate histogram');

ax[1].imshow(np.log(h[:,::-1]+1).transpose(),vmin=0,vmax=10,extent=[x.min(),x.max(),y.
˓min(),y.max()])
ax[1].set_xlabel('Input image - $f$'),ax[1].set_ylabel('$|f - \text{med}_{3x3}(f)|$'),ax[1].
˓set_title('Log bivariate histogram');
```

(continues on next page)

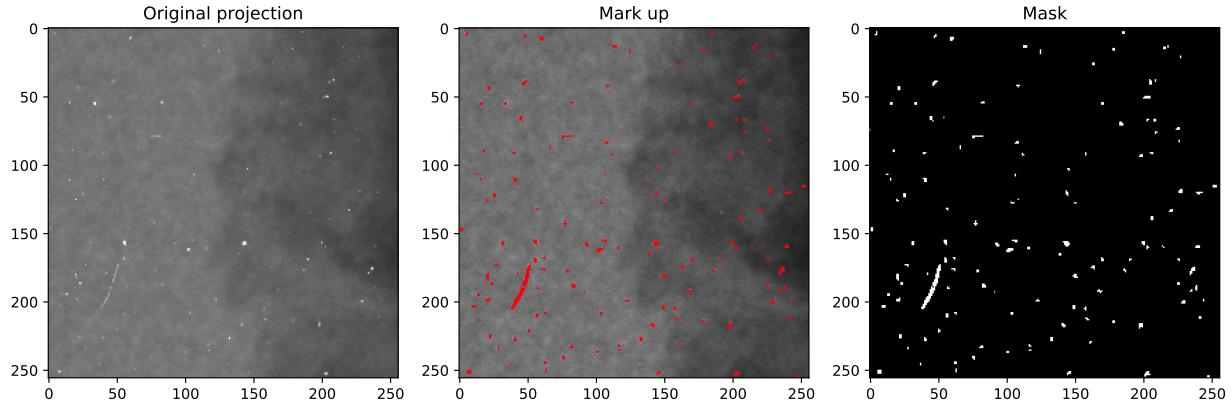


Fig. 5.3: Two cases of unbalanced data; (a) the classes are well separated and the feature class is clearly visible in the tail distribution of the background and (b) the feature class is embeded in the background making it hard to detect.

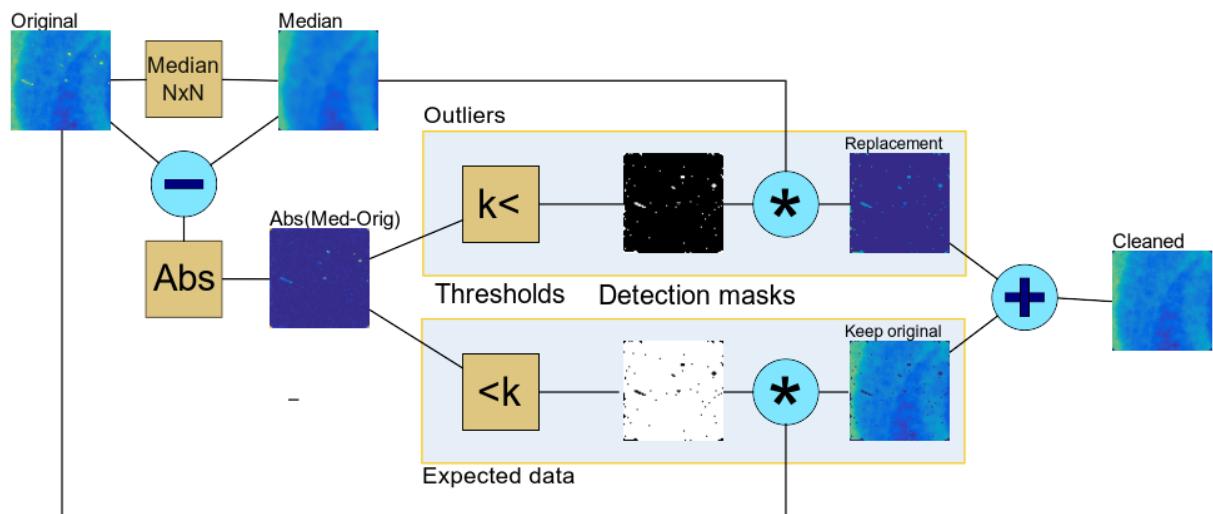
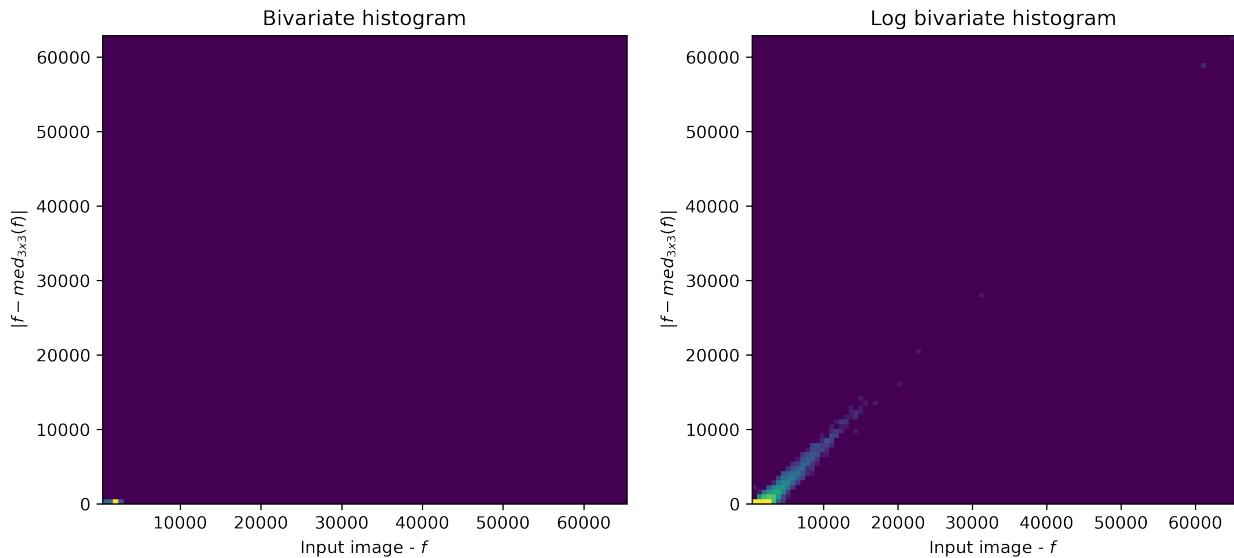


Fig. 5.4: Schematic description of the spot cleaning baseline algorithm

(continued from previous page)



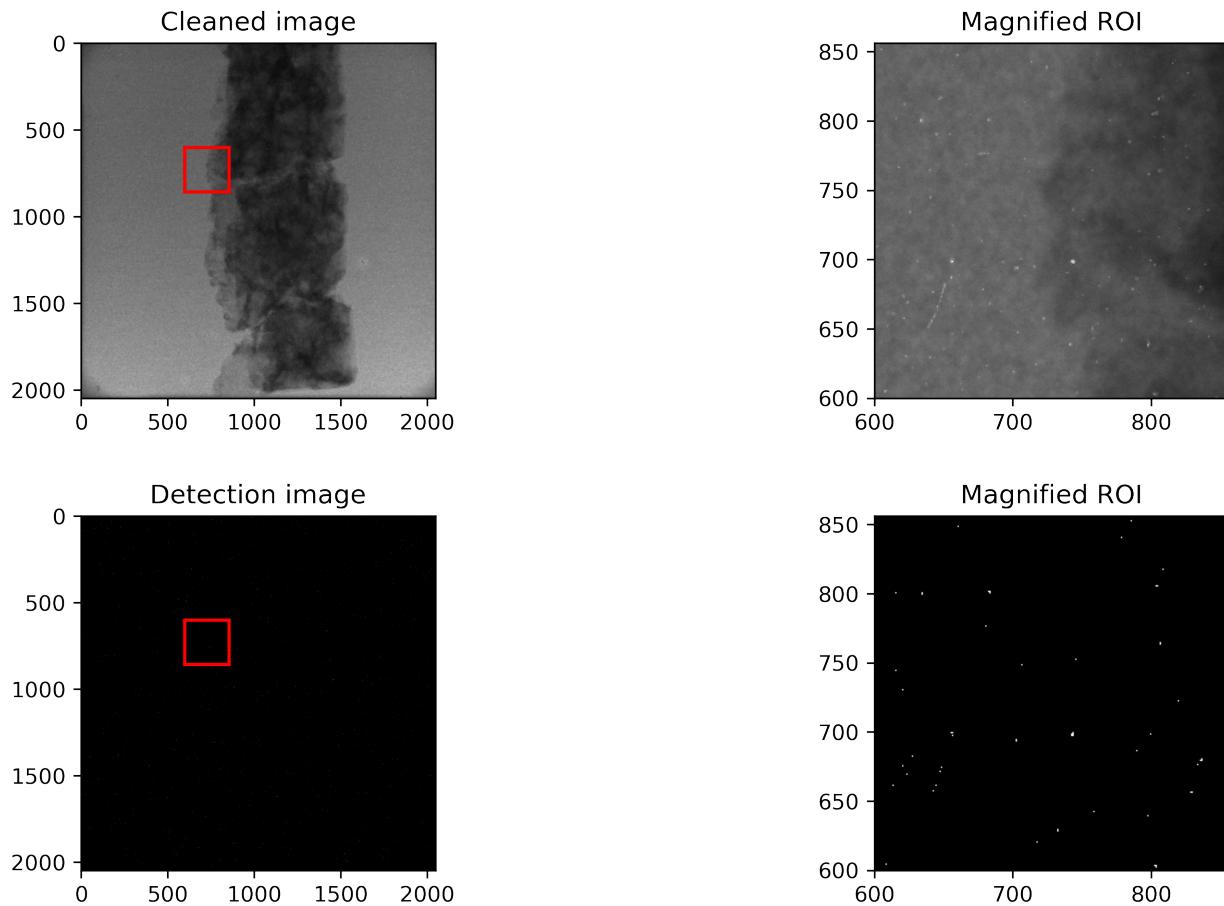
5.3.4 The spot cleaning algorithm

The baseline algorithm is here implemented as a python function that we will use when we compare the performance of the CNN algorithm. This is the most trivial algorithm for spot cleaning and there are plenty other algorithms to solve this task.

```
def spotCleaner(img, threshold=0.95, selem=np.ones([3,3])) :
    fimg=img.astype('float32')
    mimg = fimg.median(selem=selem)
    timg = threshold < np.abs(fimg-mimg)
    cleaned = mimg * timg + fimg * (1-timg)
    return (cleaned,timg)
```

5.3.5 Testing the baseline algorithm for spot cleaning

```
baseclean,timg = spotCleaner(orig,threshold=1000)
ps.magnifyRegion(baseclean,[r,c,r+w,c+w],[12,3],vmin=400,vmax=4000,title='Cleaned image')
ps.magnifyRegion(timg,[r,c,r+w,c+w],[12,3],vmin=0,vmax=1,title='Detection image')
```



5.4 k nearest neighbors to detect spots

5.4.1 Prepare data

Training data

```
trainorig = forig[:, :1000].ravel()
traind    = d[:, :1000].ravel()
trainmask = mask[:, :1000].ravel()

train_pts = pd.DataFrame({'orig': trainorig, 'd': traind, 'mask': trainmask})
```

Test data

```
testorig = forig[:, 1000: ].ravel()
testd    = d[:, 1000: ].ravel()
testmask = mask[:, 1000: ].ravel()

test_pts = pd.DataFrame({'orig': testorig, 'd': testd, 'mask': testmask})
```

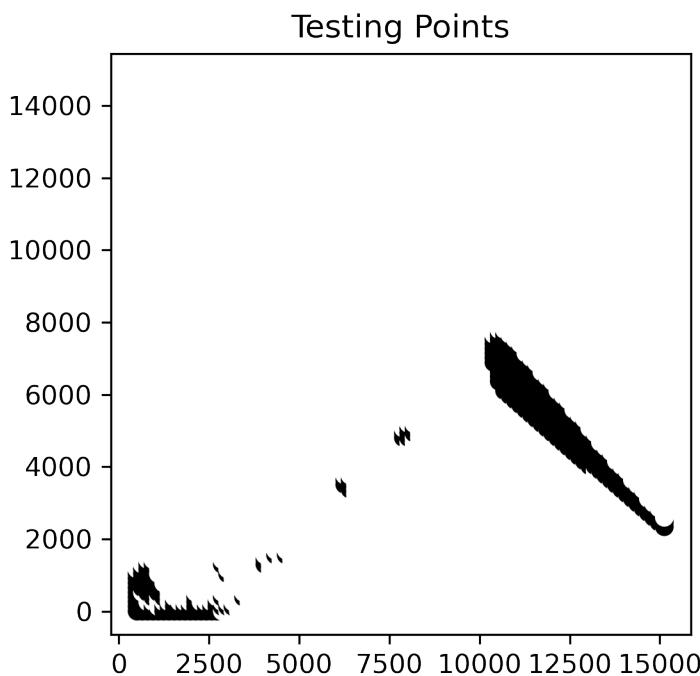
5.4.2 Train the model

```
k_class = KNeighborsClassifier(1)
k_class.fit(train_pts[['orig', 'd']], train_pts['mask'])
```

```
KNeighborsClassifier(n_neighbors=1)
```

Inspect decision space

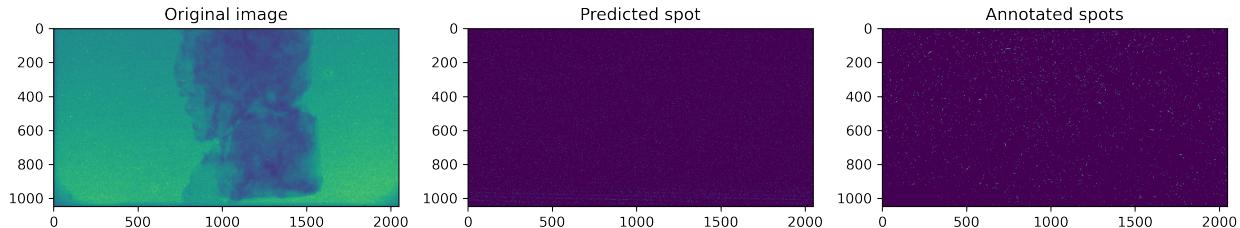
```
xx, yy = np.meshgrid(np.linspace(test_pts.orig.min(), test_pts.orig.max(), 100),
                      np.linspace(test_pts.d.min(), test_pts.d.max(), 100), indexing='ij')
grid_pts = pd.DataFrame(dict(x=xx.ravel(), y=yy.ravel()))
grid_pts['predicted_id'] = k_class.predict(grid_pts[['x', 'y']])
plt.scatter(grid_pts.x, grid_pts.y, c=grid_pts.predicted_id, cmap='gray'); plt.title(
    'Testing Points'); plt.axis('square');
```



5.4.3 Apply knn to unseen data

```
pred = k_class.predict(test_pts[['orig', 'd']])
pimg = pred.reshape(d[1000:,:].shape)
```

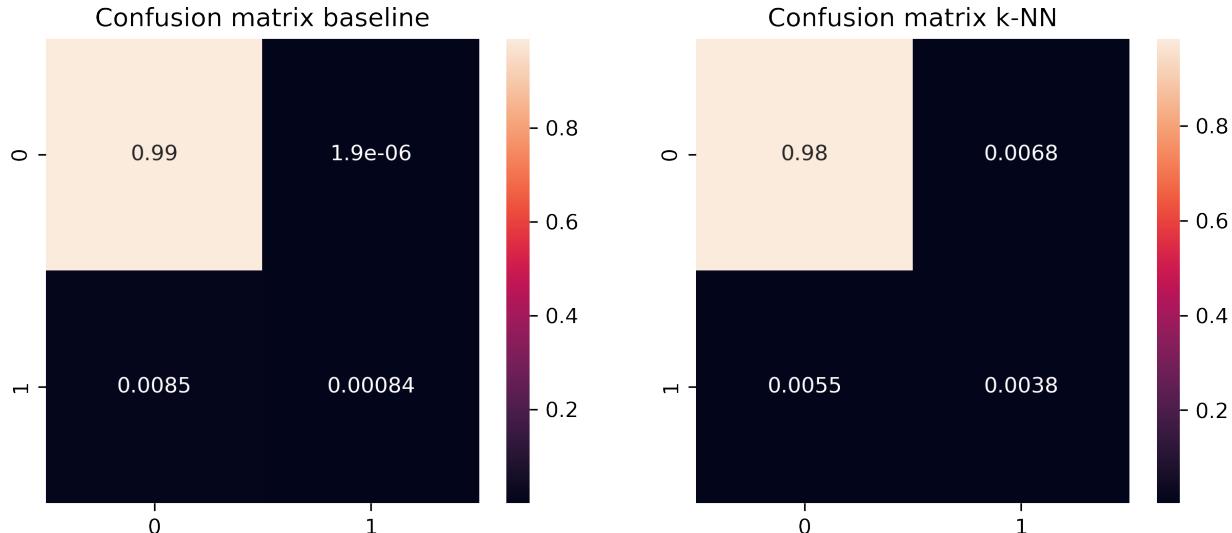
```
fig,ax = plt.subplots(1,3,figsize=(15,6))
ax[0].imshow(forig[1000,:,:],vmin=0,vmax=4000), ax[0].set_title('Original image')
ax[1].imshow(pimg), ax[1].set_title('Predicted spot')
ax[2].imshow(mask[1000,:,:]),ax[2].set_title('Annotated spots');
```



5.4.4 Performance check

```
cmbase = confusion_matrix(mask[:,1000:].ravel(), timg[:,1000:].ravel(), normalize='all')
cmknn = confusion_matrix(mask[:,1000:].ravel(), pimg.ravel(), normalize='all')
```

```
fig,ax = plt.subplots(1,2,figsize=(10,4))
sn.heatmap(cmbase, annot=True, ax=ax[0]), ax[0].set_title('Confusion matrix baseline');
sn.heatmap(cmknn, annot=True, ax=ax[1]), ax[1].set_title('Confusion matrix k-NN');
```



5.4.5 Some remarks about k-nn

- It takes more time to process
- You need to prepare training data
 - Annotation takes time...
 - Here we used the segmentation on the same type of image
 - We should normalize the data
 - This was a raw projection, what happens if we use a flat field corrected image?
- Finds more spots than baseline
- Data is very unbalanced, try a selection of non-spot data for training.
 - Is it faster?

- Is there a drop segmentation performance?

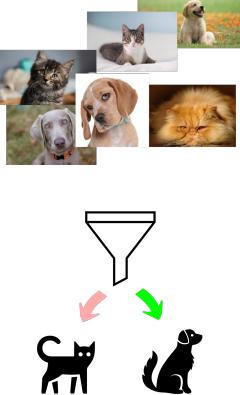
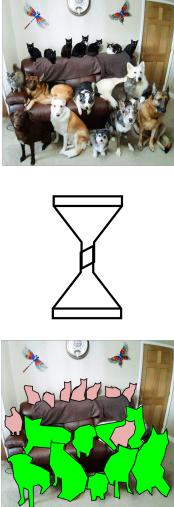
Note There are other spot detection methods that perform better than the baseline.

5.4.6 Clean up

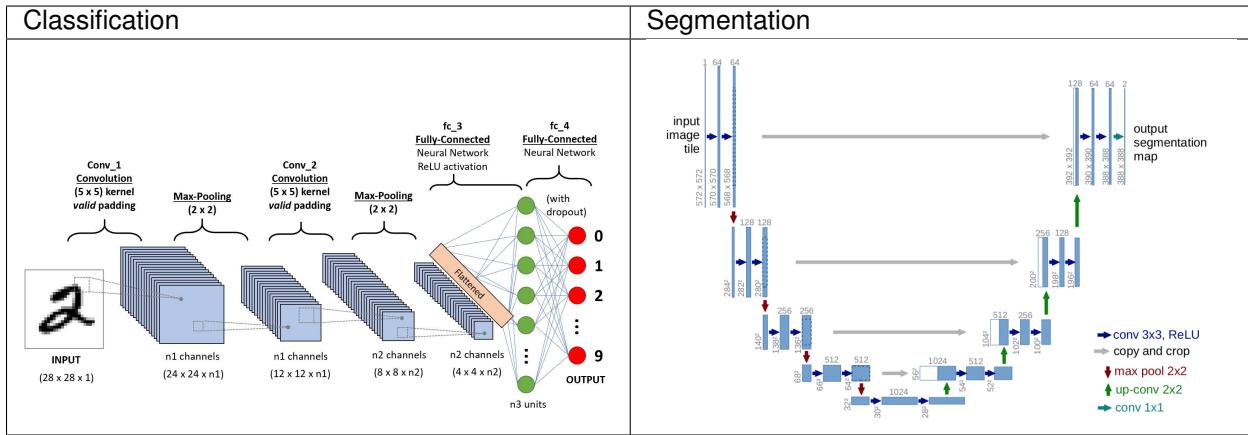
```
del k_class
```


CONVOLUTIONAL NEURAL NETWORKS FOR SEGMENTATION

6.1 The difference between classification and segmentation I

| Classification | Segmentation |
|--|---|
| pixels to classes | pixels to pixels |
|  <p>The classification section shows a grid of small images representing different animal species. Below this, a funnel icon indicates the process of reducing the complexity of the input data. Two arrows point from the funnel to two simple black silhouettes of a cat and a dog, representing the final output classes.</p> |  <p>The segmentation section shows a large image of a group of various animals. A funnel icon is shown above a second image of the same scene, but with each individual animal highlighted in a different color (pink, green, blue, etc.), representing the output of a segmentation model where each pixel is assigned to a specific object category.</p> |

6.2 Network architecture Classification vs Segmentation



6.3 Different segmentation networks

Segmentation is mostly based on variations of the U-Net architecture

- AlexNET
- SegNET
- SegCaps

6.3.1 Training data

We have two choices:

1. Use real data
 - requires time consuming markup to provide training data
 - corresponds to real life images
2. Synthesize data
 - flexible and provides both ‘dirty’ data and ground truth.
 - model may not behave as real data

6.4 Preparing real data

We will use the spotty image as training data for this example

6.4.1 Prepare training, validation, and test data

Any analysis system must be verified to be demonstrate its performance and to further optimize it.

For this we need to split our data into three categories:

1. Training data
2. Test data
3. Validation data

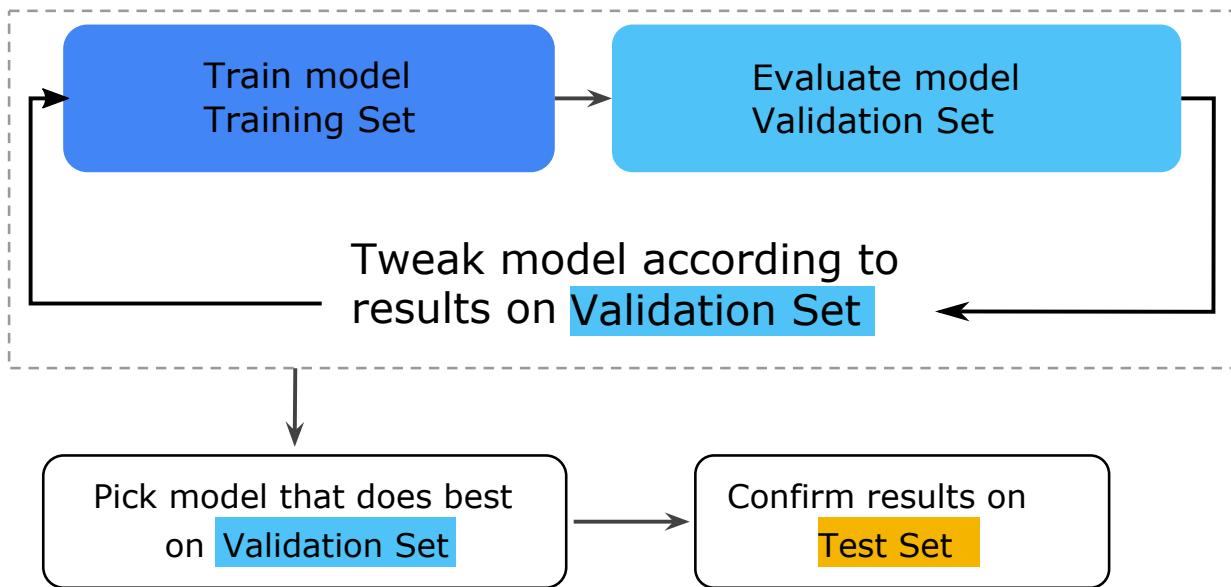


Fig. 6.1: How the three data sets *training*, *validation*, and *test* are used when a network is trained and optimized.

| Training | Validation | Test |
|----------|------------|------|
| 70% | 15% | 15% |

6.4.2 Build a CNN for spot detection and cleaning

We need:

- Data
 - Gray level image - our radiograph.
 - Annotated image where the spots are marked.
- A U-net model
 - Keras comes to our help

6.4.3 Build a U-Net model

```
def buildSpotUNet( base_depth = 48 ) :
    in_img = Input((None, None, 1), name='Image_Input')
    lay_1 = Conv2D(base_depth, kernel_size=(3, 3), padding='same',activation='relu')(in_img)
    lay_2 = Conv2D(base_depth, kernel_size=(3, 3), padding='same',activation='relu')(lay_1)
    lay_3 = MaxPooling2D(pool_size=(2, 2))(lay_2)
    lay_4 = Conv2D(base_depth*2, kernel_size=(3, 3), padding='same',activation='relu')(lay_3)
    lay_5 = Conv2D(base_depth*2, kernel_size=(3, 3), padding='same',activation='relu')(lay_4)
    lay_6 = MaxPooling2D(pool_size=(2, 2))(lay_5)
    lay_7 = Conv2D(base_depth*4, kernel_size=(3, 3), padding='same',activation='relu')(lay_6)
    lay_8 = Conv2D(base_depth*4, kernel_size=(3, 3), padding='same',activation='relu')(lay_7)
    lay_9 = UpSampling2D((2, 2))(lay_8)
    lay_10 = concatenate([lay_5, lay_9])
    lay_11 = Conv2D(base_depth*2, kernel_size=(3, 3), padding='same',activation='relu')(lay_10)
    lay_12 = Conv2D(base_depth*2, kernel_size=(3, 3), padding='same',activation='relu')(lay_11)
    lay_13 = UpSampling2D((2, 2))(lay_12)
    lay_14 = concatenate([lay_2, lay_13])
    lay_15 = Conv2D(base_depth, kernel_size=(3, 3), padding='same',activation='relu')(lay_14)
    lay_16 = Conv2D(base_depth, kernel_size=(3, 3), padding='same',activation='relu')(lay_15)
    lay_17 = Conv2D(1, kernel_size=(1, 1), padding='same',
                    activation='relu')(lay_16)
    t_unet = Model(inputs=[in_img], outputs=[lay_17], name='SpotUNET')
    return t_unet
```

Model summary

```
t_unet = buildSpotUNet(base_depth=24)
t_unet.summary()
```

WARNING:tensorflow:From /home/travis/miniconda/envs/book/lib/python3.7/site-packages/tensorflow_core/python/ops/resource_variable_ops.py:1630: calling __BaseResourceVariable.__init__ (from tensorflow.python.ops.resource_variable_ops) with constraint is deprecated and will be removed in a future version.
Instructions for updating:
If using Keras pass *_constraint arguments to layers.

WARNING:tensorflow:From /home/travis/miniconda/envs/book/lib/python3.7/site-packages/keras/backend/tensorflow_backend.py:4070: The name tf.nn.max_pool is deprecated. Please use tf.nn.max_pool2d instead.

Model: "SpotUNET"

| Layer (type) | Output Shape | Param # | Connected to |
|--------------------------|--------------|---------|--------------|
| (continues on next page) | | | |

(continued from previous page)

| | | |
|--------------------------------|-----------------------------|---|
| Image_Input (InputLayer) | (None, None, None, 1 0) | |
| conv2d_1 (Conv2D) | (None, None, None, 2 240) | Image_Input [0] [0] |
| conv2d_2 (Conv2D) | (None, None, None, 2 5208) | conv2d_1 [0] [0] |
| max_pooling2d_1 (MaxPooling2D) | (None, None, None, 2 0) | conv2d_2 [0] [0] |
| conv2d_3 (Conv2D) | (None, None, None, 4 10416) | max_pooling2d_1 [0] [0] |
| conv2d_4 (Conv2D) | (None, None, None, 4 20784) | conv2d_3 [0] [0] |
| max_pooling2d_2 (MaxPooling2D) | (None, None, None, 4 0) | conv2d_4 [0] [0] |
| conv2d_5 (Conv2D) | (None, None, None, 9 41568) | max_pooling2d_2 [0] [0] |
| conv2d_6 (Conv2D) | (None, None, None, 9 83040) | conv2d_5 [0] [0] |
| up_sampling2d_1 (UpSampling2D) | (None, None, None, 9 0) | conv2d_6 [0] [0] |
| concatenate_1 (Concatenate) | (None, None, None, 1 0) | conv2d_4 [0] [0] up_sampling2d_1 [0] [0] |
| conv2d_7 (Conv2D) | (None, None, None, 4 62256) | concatenate_1 [0] [0] |
| conv2d_8 (Conv2D) | (None, None, None, 4 20784) | conv2d_7 [0] [0] |
| up_sampling2d_2 (UpSampling2D) | (None, None, None, 4 0) | conv2d_8 [0] [0] |
| concatenate_2 (Concatenate) | (None, None, None, 7 0) | conv2d_2 [0] [0] up_sampling2d_2 [0] [0] |
| conv2d_9 (Conv2D) | (None, None, None, 2 15576) | concatenate_2 [0] [0] |
| conv2d_10 (Conv2D) | (None, None, None, 2 5208) | conv2d_9 [0] [0] |
| conv2d_11 (Conv2D) | (None, None, None, 1 25) | conv2d_10 [0] [0] |
| <hr/> | | |
| Total params: | 265,105 | |
| Trainable params: | 265,105 | |

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Segmenting neutron images using machine learning

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```
Non-trainable params: 0
```

→_____

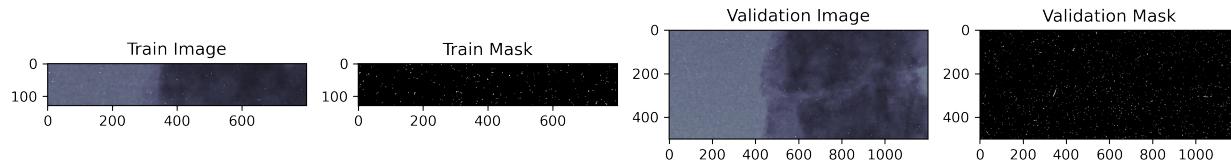
Prepare data for training and validation

```
train_img, valid_img = forig[128:256, 500:1300], forig[500:1000, 300:1500]
train_mask, valid_mask = mask[128:256, 500:1300], mask[500:1000, 300:1500]
wpos = [600, 600]; ww = 512
forigc = forig[wpos[0]:(wpos[0]+ww), wpos[1]:(wpos[1]+ww)]
maskc = mask[wpos[0]:(wpos[0]+ww), wpos[1]:(wpos[1]+ww)]

fig, ax = plt.subplots(1, 4, figsize=(15, 6), dpi=300); ax=ax.ravel()

ax[0].imshow(train_img, cmap='bone', vmin=0, vmax=4000); ax[0].set_title('Train Image')
ax[1].imshow(train_mask, cmap='bone'); ax[1].set_title('Train Mask')

ax[2].imshow(valid_img, cmap='bone', vmin=0, vmax=4000); ax[2].set_title('Validation Image')
ax[3].imshow(valid_mask, cmap='bone'); ax[3].set_title('Validation Mask');
```



Functions to prepare data for training

```
def prep_img(x, n=1):
    return (prep_mask(x, n=n)-train_img.mean()) / train_img.std()

def prep_mask(x, n=1):
    return np.stack([np.expand_dims(x, -1)]*n, 0)
```

Test the untrained model

- We can make predictions with an untrained model (default parameters)
- but we clearly do not expect them to be very good

```
unet_pred = t_unet.predict(prep_img(forigc))[0, :, :, 0]
```

WARNING:tensorflow:From /home/travis/miniconda/envs/book/lib/python3.7/site-packages/keras/backend/tensorflow_backend.py:422: The name tf.global_variables is deprecated.
→ Please use tf.compat.v1.global_variables instead.

```
fig, m_axs = plt.subplots(2, 3, figsize=(15, 6), dpi=150)
for c_ax in m_axs.ravel():
```

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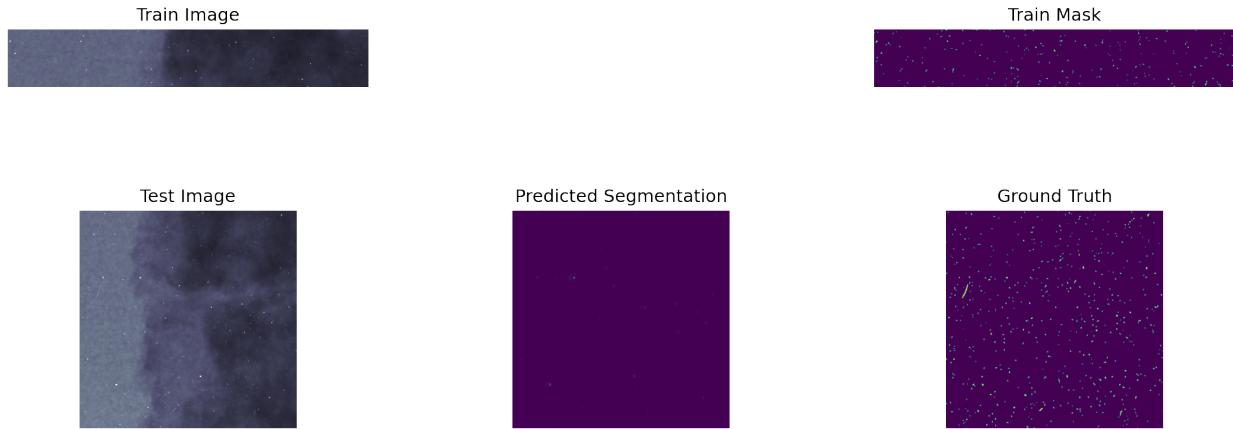
(continued from previous page)

```

c_ax.axis('off')
((ax1, _, ax2), (ax3, ax4, ax5)) = m_axs
ax1.imshow(train_img, cmap='bone', vmin=0, vmax=4000); ax1.set_title('Train Image')
ax2.imshow(train_mask, cmap='viridis'); ax2.set_title('Train Mask')

ax3.imshow(forigc, cmap='bone', vmin=0, vmax=4000); ax3.set_title('Test Image')
ax4.imshow(unet_pred, cmap='viridis', vmin=0, vmax=1); ax4.set_title('Predicted  
Segmentation')

ax5.imshow(maskc, cmap='viridis'); ax5.set_title('Ground Truth');
    
```



6.4.4 Training conditions

- Loss function - Binary cross-correlation
- Optimizer - ADAM
- 20 Epochs (training iterations)
- Metrics
 1. True positives
 2. False positives
 3. True negatives
 4. False negatives
 5. Binary accuracy (percentage of pixels correctly classified) $BA = \frac{1}{N} \sum_i (f_i == g_i)$
 6. Precision $Precision = \frac{TP}{TP+FP}$
 7. Recall $Recall = \frac{TP}{TP+FN}$
 8. Area under receiver operating characteristics (ROC) curve $AUC = \int ROC$
 9. Mean absolute error $MAE = \frac{1}{N} \sum_i |f_i - g_i|$

6.5 Compile the model

```
mlist = [
    metrics.TruePositives(name='tp'),           metrics.FalsePositives(name='fp'),
    metrics.TrueNegatives(name='tn'),           metrics.FalseNegatives(name='fn'),
    metrics.BinaryAccuracy(name='accuracy'),   metrics.Precision(name='precision'),
    metrics.Recall(name='recall'),             metrics.AUC(name='auc'),
    metrics.MeanAbsoluteError(name='mae'))]

t_unet.compile(
    loss=loss.BinaryCrossentropy(),      # we use the binary cross-entropy to optimize
    optimizer=opt.Adam(lr=1e-3),         # we use ADAM to optimize
    metrics=mlist                        # we keep track of the metrics in mlist
)
```

```
WARNING:tensorflow:From /home/travis/miniconda/envs/book/lib/python3.7/site-packages/
        ↪keras/backend/tensorflow_backend.py:3172: where (from tensorflow.python.ops.array_
        ↪ops) is deprecated and will be removed in a future version.
Instructions for updating:
Use tf.where in 2.0, which has the same broadcast rule as np.where
```

6.5.1 A general note on the following demo

This is a very bad way to train a model;

- the loss function is poorly chosen,
- the optimizer can be improved the learning rate can be changed,
- the training and validation data **should not** come from the same sample (and **definitely** not the same measurement).

The goal is to be aware of these techniques and have a feeling for how they can work for complex problems

6.5.2 Training the spot detection model

```
loss_history = t_unet.fit(prep_img(train_img, n=3),
                           prep_mask(train_mask, n=3),
                           validation_data=(prep_img(valid_img),
                                             prep_mask(valid_mask)),
                           epochs=20,
                           verbose = 0)
```

Training history plots

```
titleDict = {'tp': "True Positives", 'fp': "False Positives", 'tn': "True Negatives", 'fn':
        ↪": "False Negatives", 'accuracy': "BinaryAccuracy", 'precision': "Precision", 'recall':
        ↪": "Recall", 'auc': "Area under Curve", 'mae': "Mean absolute error"}

fig,ax = plt.subplots(2,5, figsize=(20,8), dpi=300)
ax=ax.ravel()
for idx,key in enumerate(titleDict.keys()):
    ax[idx].plot(loss_history.epoch, loss_history.history[key], color='coral', label=
        ↪": "Training")
```

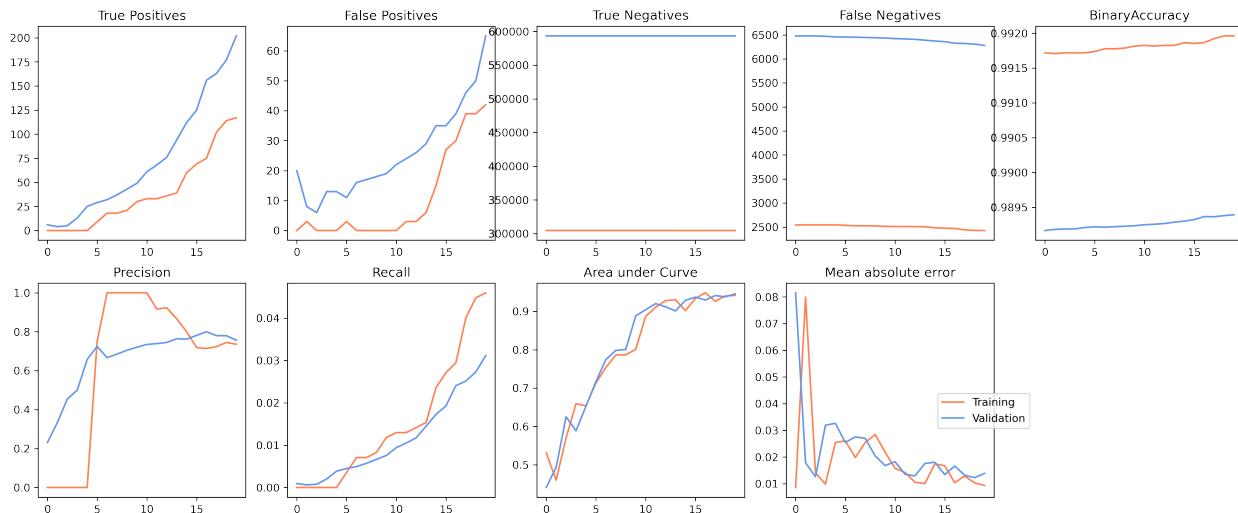
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```

        ax[idx].plot(loss_history.epoch, loss_history.history['val_'+key], color=
        ↪'cornflowerblue', label='Validation')
        ax[idx].set_title(titleDict[key]);

ax[9].axis('off');
axLine, axLabel = ax[0].get_legend_handles_labels() # Take the labels and plot line
↪information from the first panel
lines =[]; labels = []; lines.extend(axLine); labels.extend(axLabel);fig.legend(lines,
↪ labels, bbox_to_anchor=(0.7, 0.3), loc='upper left');
    
```



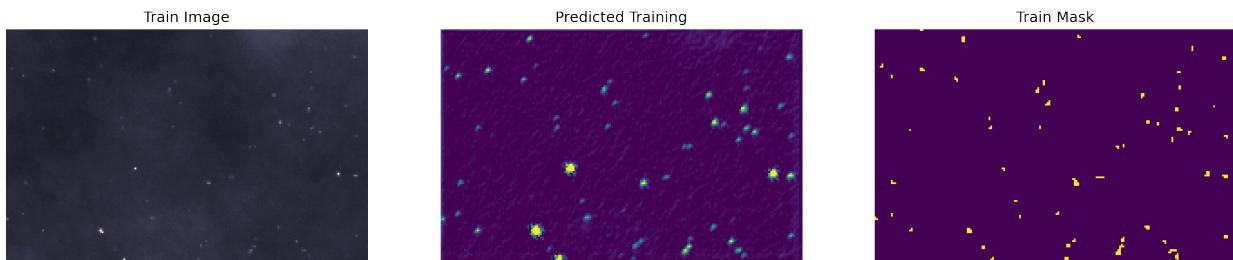
Prediction on the training data

```

unet_train_pred = t_unet.predict(prep_img(train_img[:,wpos[1]:(wpos[1]+ww)])) [0, :, :, 
↪ 0]

fig, m_axs = plt.subplots(1, 3, figsize=(18, 4), dpi=150); m_axs= m_axs.ravel();
for c_ax in m_axs: c_ax.axis('off')

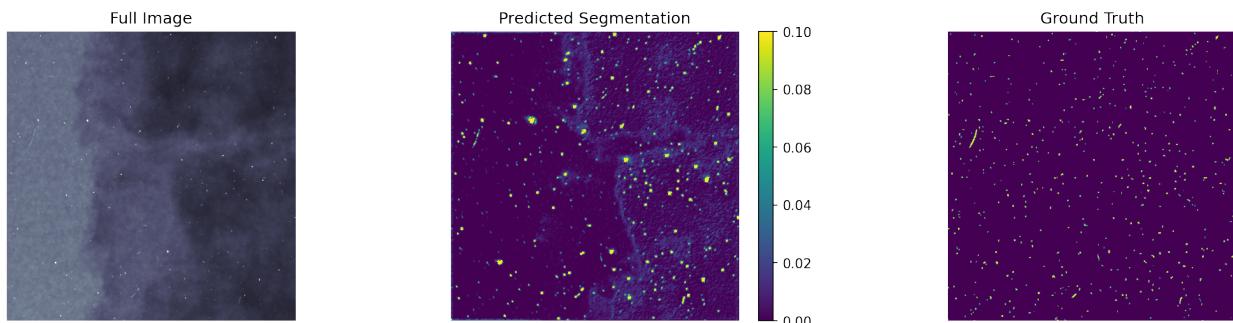
m_axs[0].imshow(train_img[:,wpos[1]:(wpos[1]+ww)], cmap='bone', vmin=0, vmax=4000), m_
↪axs[0].set_title('Train Image')
m_axs[1].imshow(unet_train_pred, cmap='viridis', vmin=0, vmax=0.2), m_axs[1].set_
↪title('Predicted Training')
m_axs[2].imshow(train_mask[:,wpos[1]:(wpos[1]+ww)], cmap='viridis'), m_axs[2].set_
↪title('Train Mask');
    
```



6.5.3 Prediction using unseen data

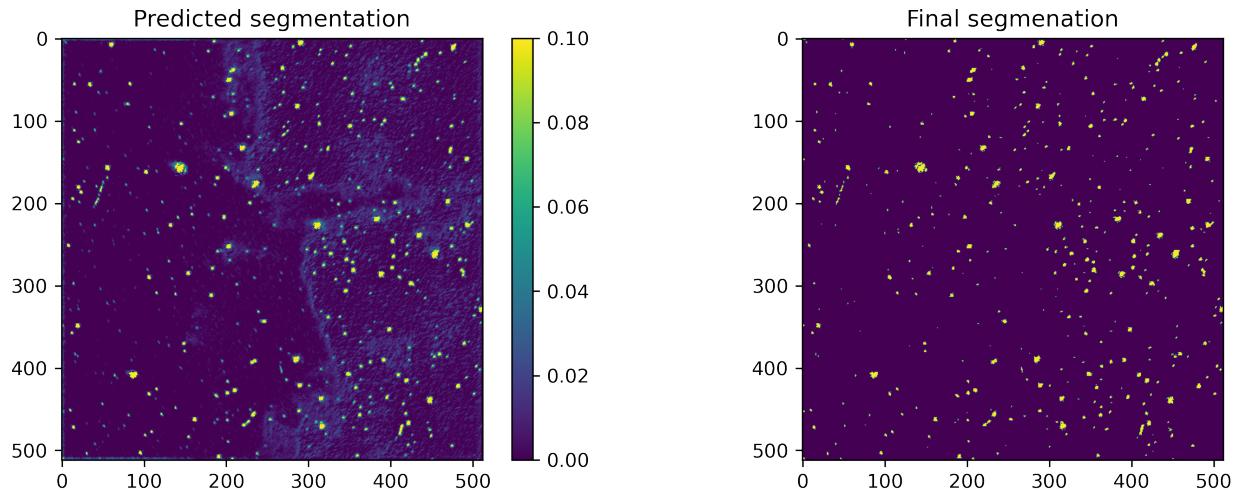
```
unet_pred = t_unet.predict(prep_img(forigc)) [0, :, :, 0]

fig, m_axs = plt.subplots(1, 3, figsize=(18, 4), dpi=150); m_axs = m_axs.ravel() ;
for c_ax in m_axs: c_ax.axis('off')
m_axs[0].imshow(forigc, cmap='bone', vmin=0, vmax=4000); m_axs[0].set_title('Full Image')
f1=m_axs[1].imshow(unet_pred, cmap='viridis', vmin=0, vmax=0.1); m_axs[1].set_title('Predicted Segmentation')
m_axs[2].imshow(maskc,cmap='viridis'); m_axs[2].set_title('Ground Truth');
```



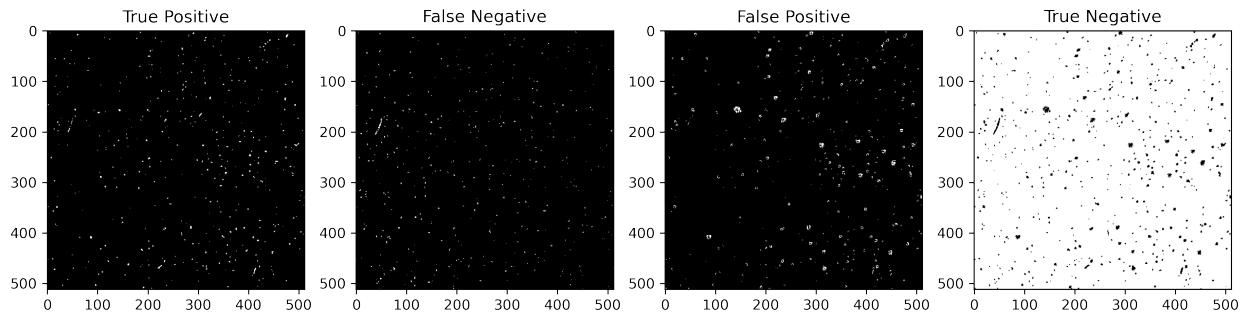
Converting predictions to segments

```
fig, ax = plt.subplots(1,2, figsize=(12,4))
ax0=ax[0].imshow(unet_pred, vmin=0, vmax=0.1); ax[0].set_title('Predicted segmentation')
fig.colorbar(ax0,ax=ax[0])
ax[1].imshow(0.05<unet_pred), ax[1].set_title('Final segmenation');
```



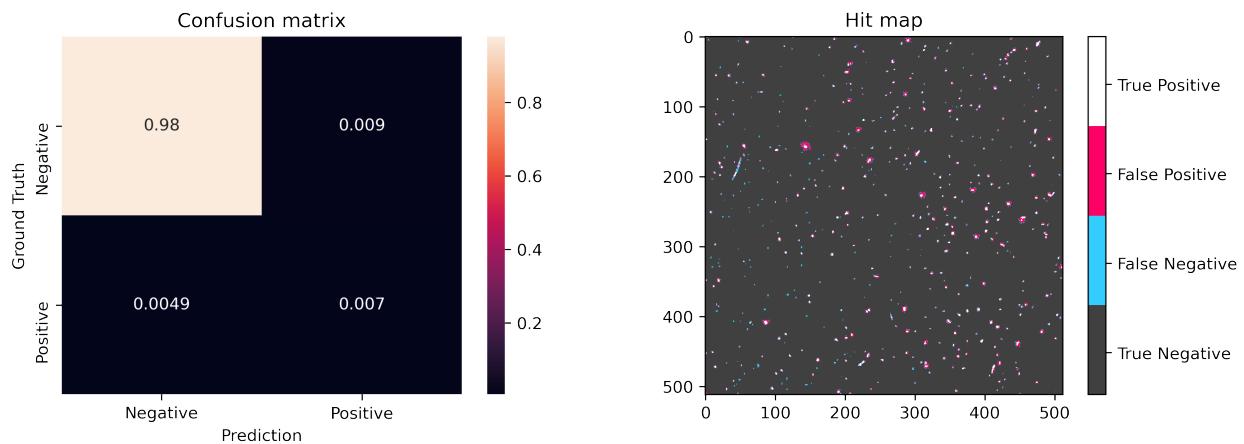
Hit cases

```
gt = maskc
pr = 0.05<unet_pred
ps.showHitCases(gt,pr,cmap='gray')
```



Hit map

```
fig, ax = plt.subplots(1,2,figsize=(12,4))
ps.showHitMap(gt,pr,ax=ax)
```

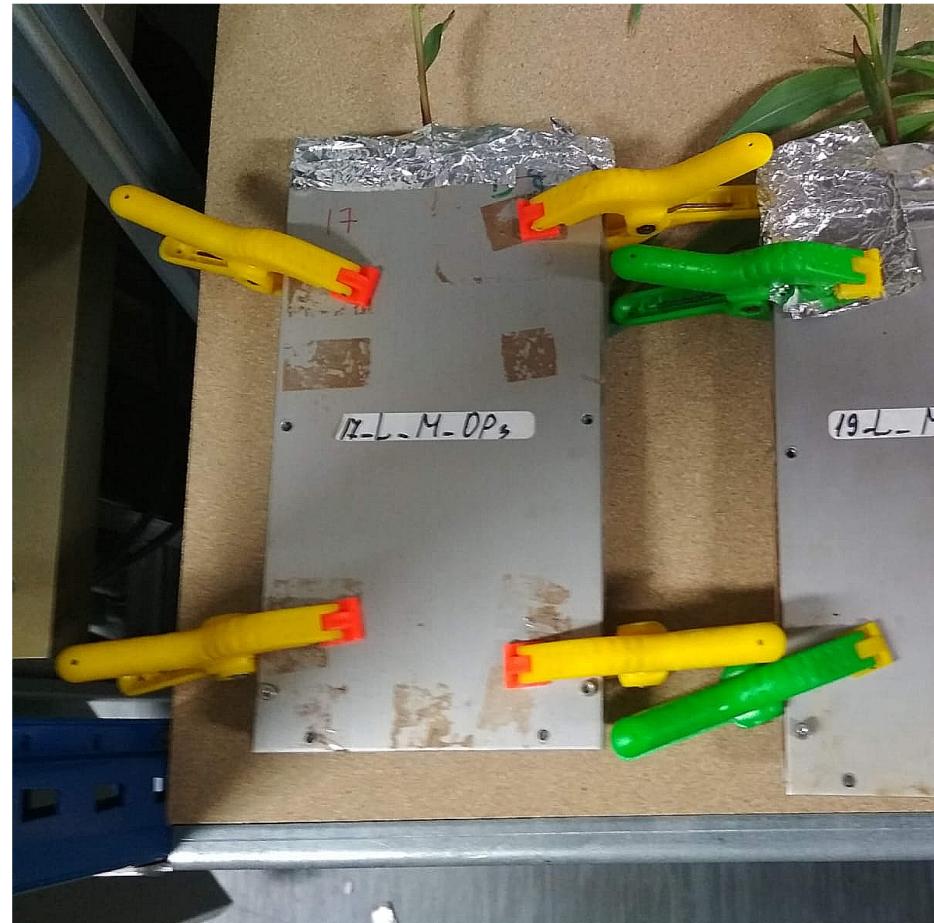


6.5.4 Concluding remarks about the spot detection

SEGMENTING ROOT NETWORKS IN THE RHIZOSPHERE

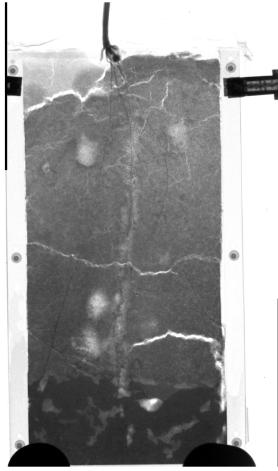
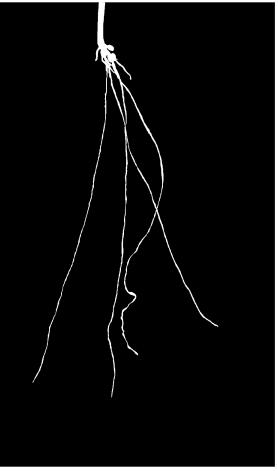
7.1 Background

- Soil and in particular the rhizosphere are of central interest for neutron imaging users.
- The experiments aim to follow the water distribution near the roots.
- The roots must be identified in 2D and 3D data



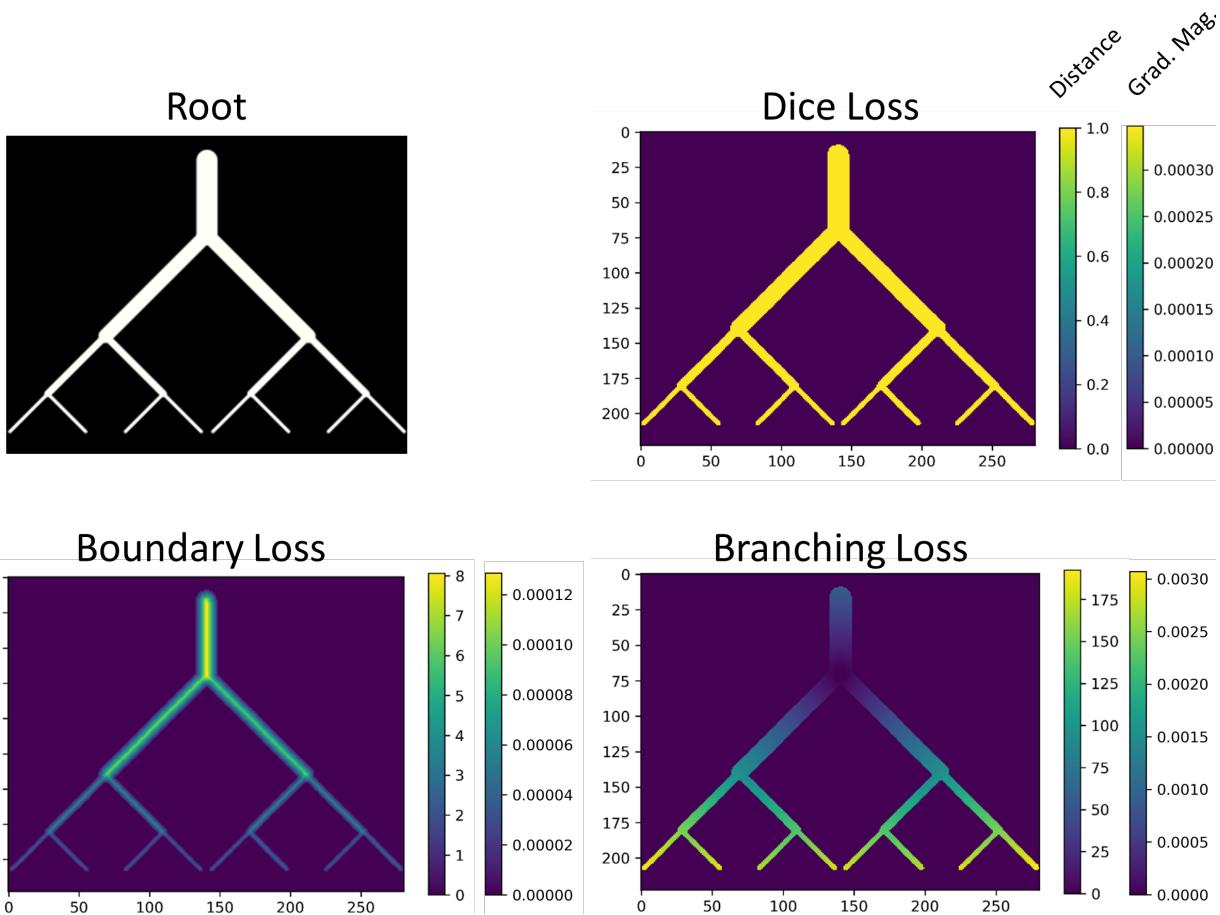
Today: much of this mark-up is done manually!

7.2 Available data

| Radiography | Tomography |
|---|---|
|  |  |
| Provided by A. Carminati et al. | Provided by M. Menon et al. |

7.3 Considered NN models

7.4 Loss functions



7.5 Training

| | Radiography | Tomography |
|-----------|-------------|------------|
| Data size | 256x256 | 64x64x64 |

7.6 Results

7.7 Summary

**CHAPTER
EIGHT**

CONCLUDING REMARKS

We have demonstrated how some machine learning techniques can be used on neutron images:

- Some background to the segmentation problem and neutron imaging
- k-means - to segment ToF spectra
- Spot detection
 - Baseline algorithm
 - k-Nearest neighbors
 - U-Net
- Root segmentation
 - U-Net
 - Different loss metrics
 - Training performance