
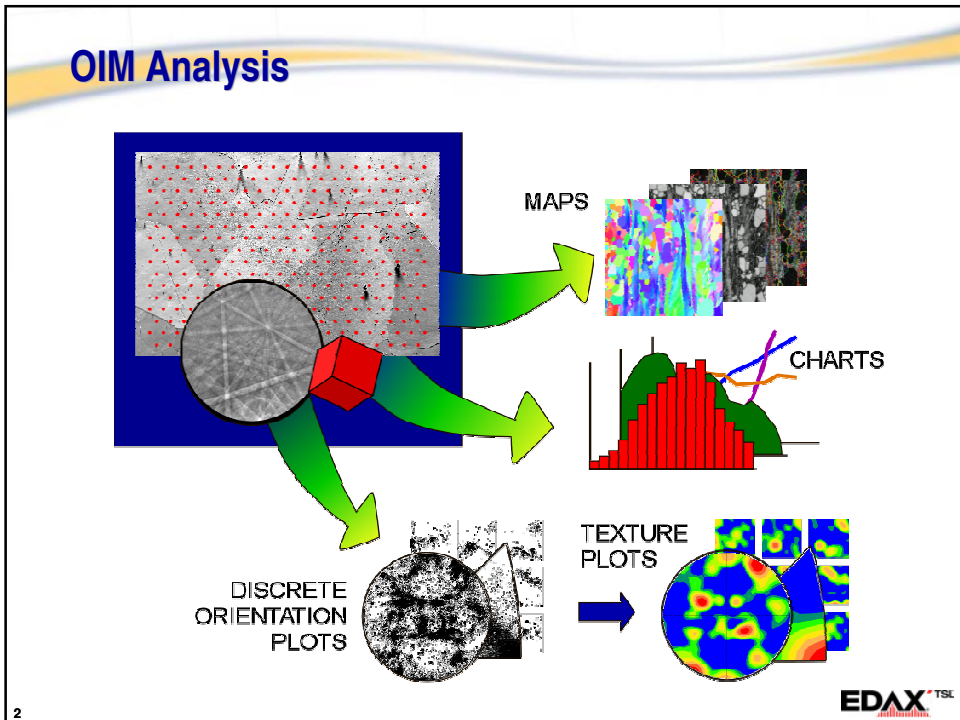


Introduction to OIM Analysis



EDAX^{TSL}
advanced microanalysis solutions
AMETEK

This slide features a background with a grid and a blue-to-orange gradient wave. On the left, there is an illustration of a scanning electron microscope (SEM) with an energy-dispersive X-ray (EDX) detector. The title 'Introduction to OIM Analysis' is centered in a large blue font.




OIM Analysis

MAPS

CHARTS

DISCRETE ORIENTATION PLOTS

TEXTURE PLOTS



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The diagram illustrates the OIM analysis workflow. It starts with a central image of a grain structure with red dots representing dislocations. Three green arrows point from this central image to three different analysis outputs: 1) 'MAPS' (Orientation Mapping) showing a color-coded grain orientation map; 2) 'CHARTS' showing a histogram of grain orientations with a red bar chart and a purple line graph; 3) 'DISCRETE ORIENTATION PLOTS' showing a circular plot of discrete orientations. A blue arrow points from the 'DISCRETE ORIENTATION PLOTS' to 'TEXTURE PLOTS', which shows a circular texture plot with a color gradient representing orientation density.

OIM Analysis

The screenshot shows the OIM Analysis software interface. At the top, there are three toolbars: **Highlighting Toolbar**, **Standard Toolbar**, and **Crystal Utilities Toolbar**. On the left side, there is a **Project Tree** and a **QuickGen Toolbar**. The main workspace is divided into several panes: a **Chart** (Pole Plot) showing a line graph of Intensity vs. Angle, a **Map** showing a color-coded orientation map, a **Summary Window** displaying statistical data like 'Name: All data', 'Formula: $P(\theta) \propto 0.100$ ', and 'Number of points in partition: 37298', and a **Texture Plot** showing a circular orientation distribution. At the bottom, there is a **Status bar** and a **Document Selector**. The **Main View Pane** and **Secondary View Pane** are also indicated. The EDAX logo is visible in the bottom right corner.

Project tree

The screenshot shows the project tree in OIM Analysis 4.5. The tree structure is as follows:

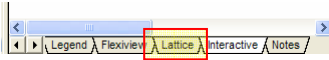
- Sample Project** (Project)
 - Grain Size Multichart (A chart that combines results from any partition in the project)
 - Fiber cleaned (OIM Datasets)
 - All data (OIM Datasets)
 - New GB Texture (Partitions. Essentially subsets of the data that meet some user prescribed criteria.)
 - MDF (Partitions. Essentially subsets of the data that meet some user prescribed criteria.)
 - GB Multi-chart (A "Chart" – e.g. grain size distribution, IQ distribution...)
 - Fiber (A map)
 - H CI data (A discrete orientation plot such as a pole figure or an "ODF" – i.e. discrete points plotted in Euler space.)
 - Auto GS (diameter) (A "Texture" – this contains all of the results for a given calculations. i.e. for a user prescribed set of parameters.)
 - Auto IPF (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - Auto IPF (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - Lo CI data (A discrete orientation plot such as a pole figure or an "ODF" – i.e. discrete points plotted in Euler space.)
 - Auto GS (diameter) (A "Texture" – this contains all of the results for a given calculations. i.e. for a user prescribed set of parameters.)
 - Auto IPF (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - Auto IPF (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - 60% Res (A discrete orientation plot such as a pole figure or an "ODF" – i.e. discrete points plotted in Euler space.)
 - All data (A "Texture" – this contains all of the results for a given calculations. i.e. for a user prescribed set of parameters.)
 - Harmonic: L=16, HW=5.0 (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - ODF (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - IPF (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - IPF (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - Euler ODF (Texture plots (of intensities) such as a Pole Figures or ODFs.)
 - IQ + Orientation Deviation (Texture plots (of intensities) such as a Pole Figures or ODFs.)

Practice session – OIM Analysis

- Create a project
 - File Menu or use icon on toolbar



- Right mouse click on project icon in project tree and select New>Dataset.
- Open “Dual Phase Titanium.osc”
- Right mouse click on “All Data” partition in the project tree and select New>Map
- Press the “OK” button in the Map Properties Dialog
- Switch to the “Lattice” pane in the Map Window

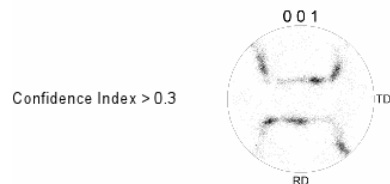
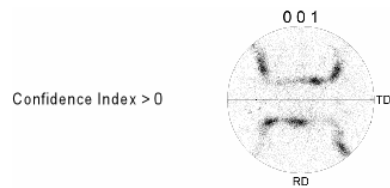
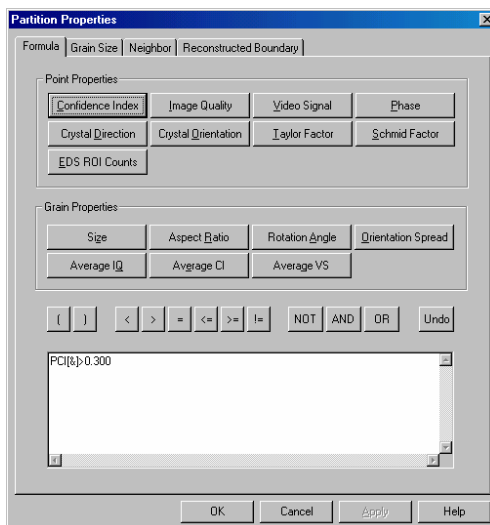


- Move the cursor around the map and note the rotating crystal
- Note the structure of the project tree

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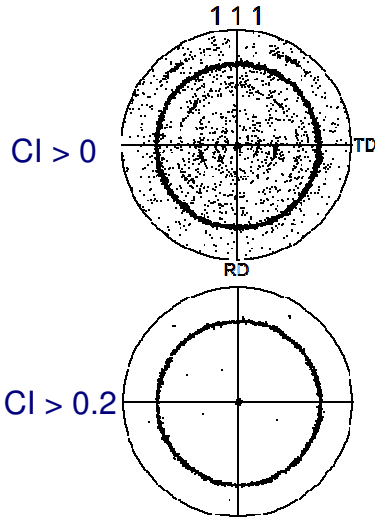
Partitions



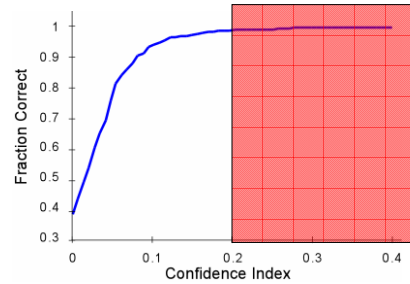
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Partitions



The data can be filtered based on any variety of parameters. In this case the confidence index.



Aluminum Thin Film

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“Grains” in OIM

While the concept of a grain in conventional metallography is commonly understood, the definition of a grain in an OIM scan is slightly different .

Grains in OIM are formed by an algorithm that groups sets of connected and similarly oriented points into “grains”.

For each point in the OIM scan, the neighbors of this point are checked to see if they are within the Grain Tolerance Angle of the given point.

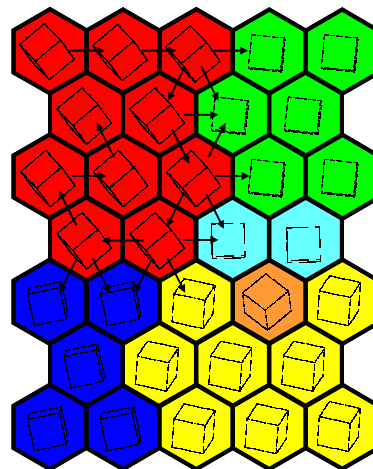
If a neighboring point is found to be within the tolerance angle then the neighbors of this point are checked to see if they are within the tolerance angle of this point.

The procedure is repeated over and over again until the set of connected grains is bounded by points which exceed the tolerance angle.

Using this approach, the point to point misorientation in a “grain” will be quite small but the spread of orientation among all points in the “grain” can be relatively large.

The number of points required to decide whether a given group of points should be considered a “grain” group (the Minimum Grain Size) can be specified by the user along with the Grain Tolerance Angle.

Thus, the definition of a grain in OIM can vary depending on user-specified values. These parameters are set using the Grain tab in the partition dialog.

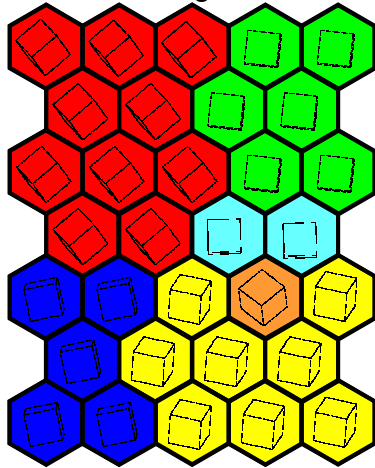


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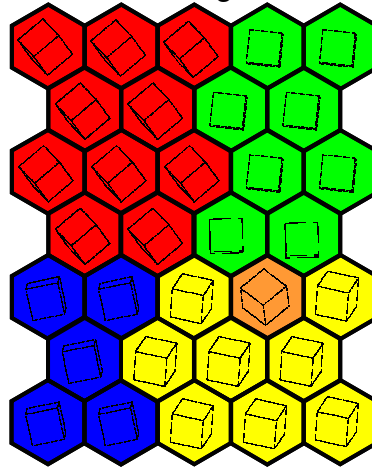
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Grain size parameters – tolerance angle

1 degree



15 degrees

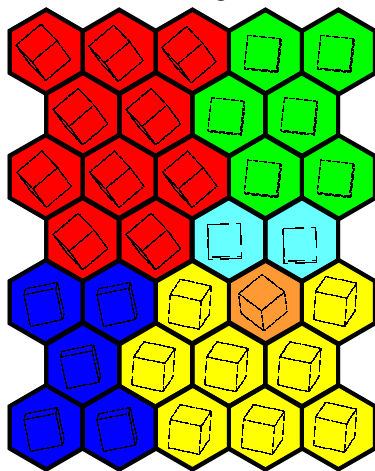


9

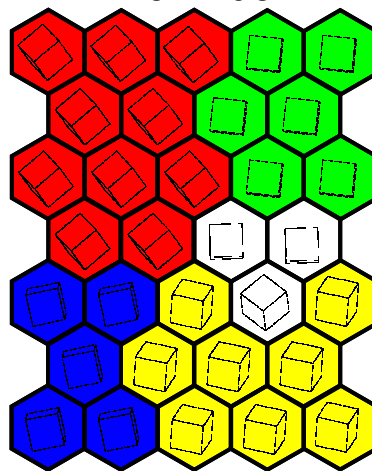
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Grain size parameters – min grain size

1 "Pixel"



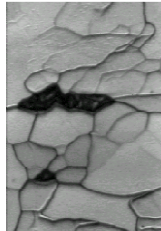
3 "Pixels"



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“Grains” in OIM

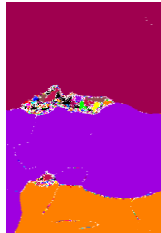


Effect of Grain Tolerance Angle



2.5µm = 50 steps

10 degrees



1 degrees



0.5 degrees

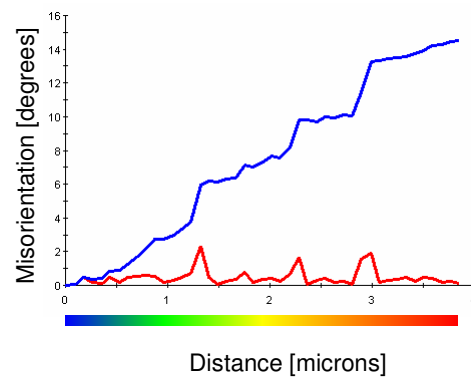
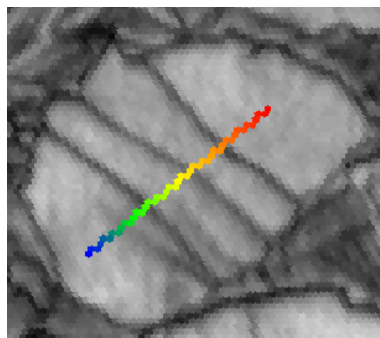


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Grain size - misorientations

The grain grouping algorithm focuses only on neighbor-to-neighbor misorientations. Thus, while the pixel-to-pixel misorientations will be less than the tolerance angle; the misorientation from one side of the grain to the other can be quite large. This is particularly evident in deformed materials.



— Point-to-point misorientation
— Point-to-origin misorientation

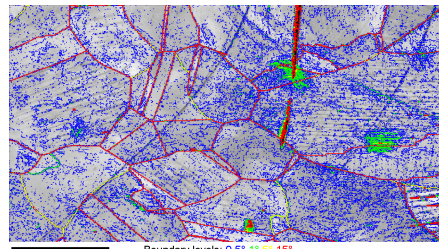
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“Grains” in OIM

Choosing a cut-off value is not always obvious – in fact, the choice of cutoff value is often best selected by visual inspection. In the case shown here of choice of 5 degrees seems quite reasonable.

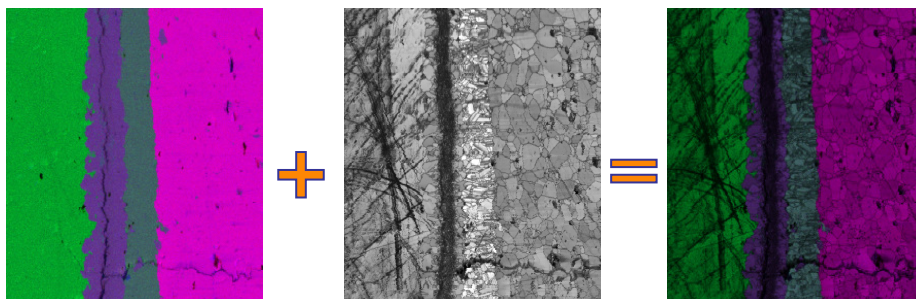
OIM gives a more precise meaning to the term boundary. However, there is obviously a great deal of literature in materials science that uses the definition from conventional metallography. It is not always an easy comparison between OIM and conventional metallography.



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Maps



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Map types

- ➔ Orientation based maps
 - Inverse Pole Figure
 - Crystal Direction
 - Crystal Orientation
- ➔ Misorientation based maps
 - Kernel Average Misorientation
 - Grain Orientation Spread
 - Grain Average Misorientation
 - Grain Reference Orientation Deviation
 - Local Average Misorientation
 - Local Orientation Spread
- ➔ Non-orientation scalar maps
 - Image Quality
 - Confidence Index
 - Fit
 - Detector Signal
 - Phase
 - EDS
 - EDS Blended Color
- ➔ Grain based maps
 - Unique Grain Color
 - Grain Size
 - Grain Shape Orientation
 - Grain Shape Aspect Ratio
 - Grain Shape Major Axis
 - Grain Shape Minor Axis
 - Major Axis Inverse Pole Figure
 - Grain Average IQ
 - Grain Average CI
 - Grain Average Fit
 - Grain Average Video Signal
 - Twin Parent/Daughter
- ➔ Property maps
 - Taylor Factor
 - Schmid Factor
 - Elastic Stiffness
- ➔ Custom
 - Import Data

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OIM: IQ, CI, video signal

Image Quality

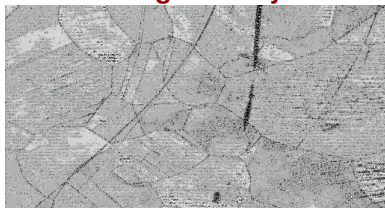
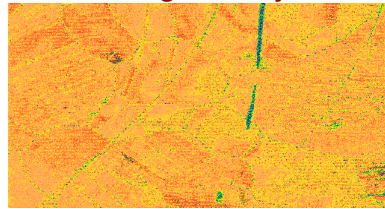
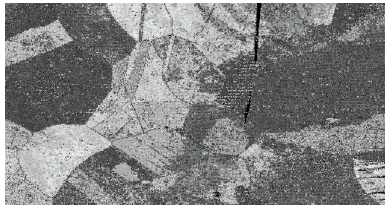


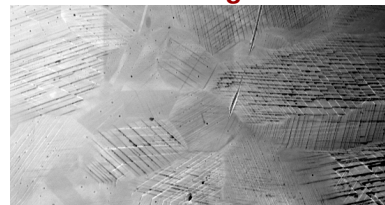
Image Quality



Confidence index



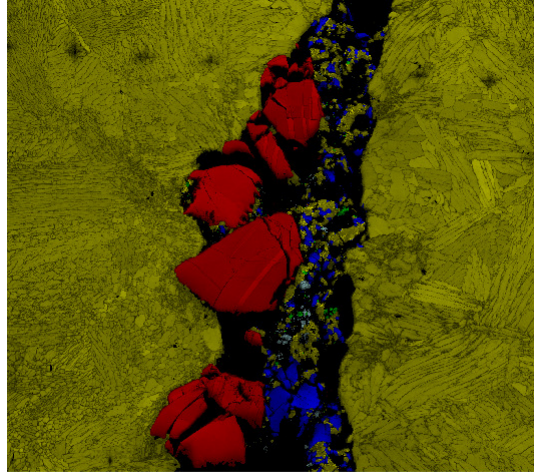
Video signal



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Maps: phase



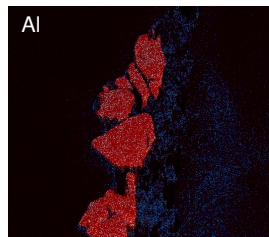
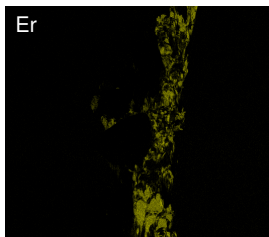
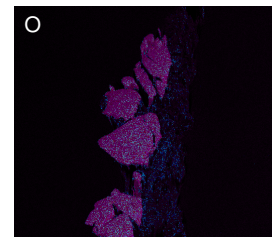
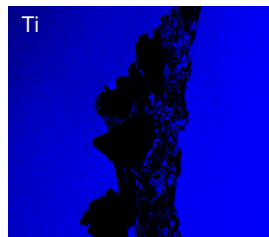
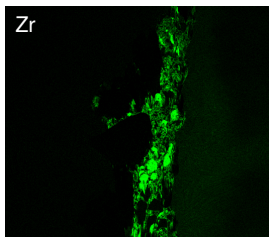
- Aluminum Titanium
- Alumina
- Erbium Oxide
- Zirconium Oxide

100.0 μm = 100 steps IQ 0...51.284, Phase

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Maps: EDS

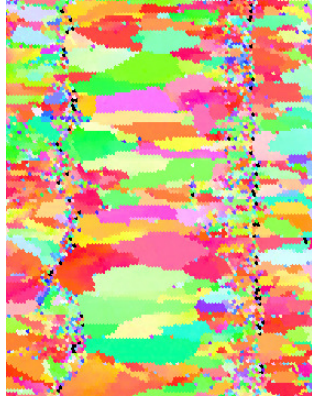


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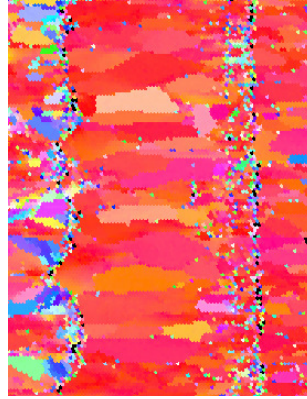
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Maps: orientation based

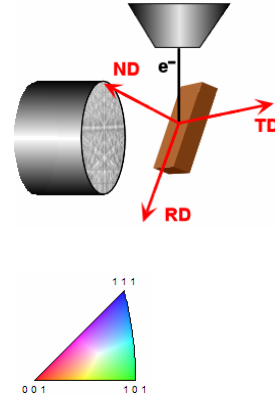
IPF maps (2D)



Normal direction



Transverse direction

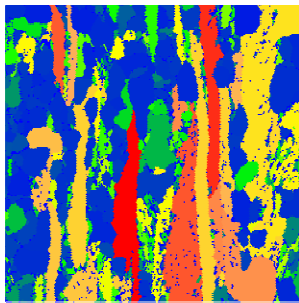


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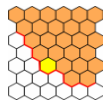
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OIM: local misorientation

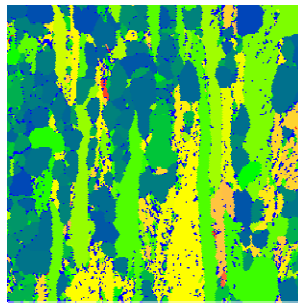
Grain Orientation Spread
For a given grain calculate the average misorientation between all data points in the grain.



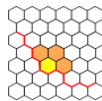
33.70 μm = 40 steps Grain Spread G...14.3603



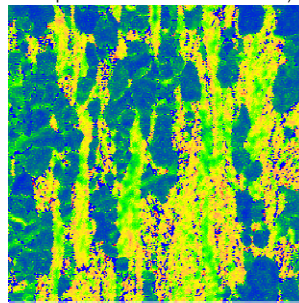
Grain Average Misorientation
For a given grain calculate the average misorientation between all neighboring data points in the grain.



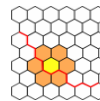
33.70 μm = 40 steps Grain Average Misorientation G...4.04363



Kernel Average Misorientation
For a given data point calculate the average misorientation between the data point and all of its neighbors (exclude misorientations greater than some prescribed value - 5° in this case)



33.70 μm = 40 steps Kernel Misorientation G...4.36361



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Practice session - mapping

- ➔ Open CVDTa.osc
- ➔ Create an IQ map using the QuickGen button



- ➔ Do a right mouse click on the map and select properties
- ➔ Set "Color Coded" to crystal direction
- ➔ Press the "Edit >>" Button
- ➔ Press the "Add" Button
- ➔ Set Sample Direction to RD=0, TD=1 & ND=0
- ➔ Switch to Color Gradient and choose a color scheme
- ➔ Look at the map legend – does it make sense?
 - Try changing the tolerance value and note effect

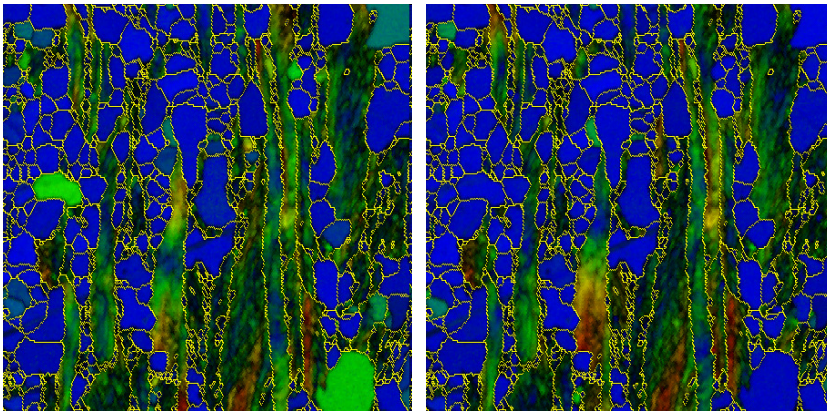
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Maps: local misorientation

Grain Reference Orientation Deviation

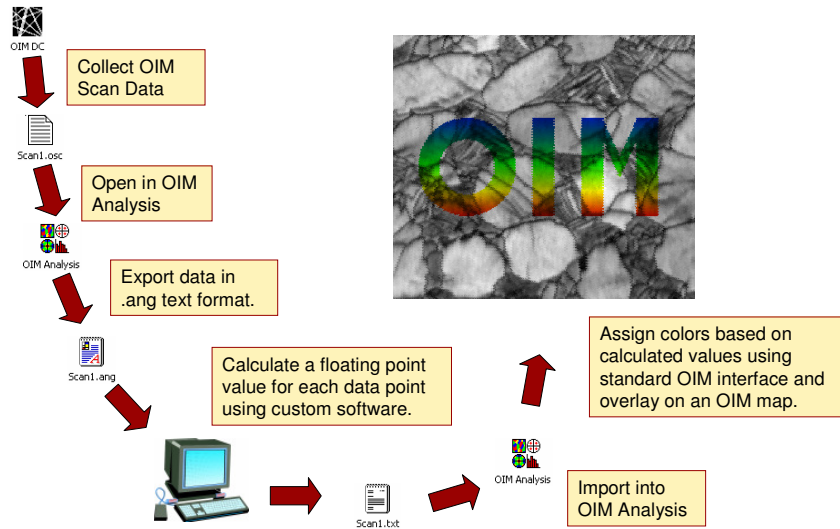
Show the deviation of each point in the grain with respect to a reference orientation. Two types of references are possible: 1) The average orientation for the grain and 2) the orientation of the point in the grain with the lowest Average Kernel Misorientation.



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Maps: import



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Boundaries

Rotation angle – A line segment is drawn between two neighboring points if the misorientation falls in a given range (e.g. between 5 and 15 degrees)

Rotation axis – A line segment is drawn between two neighboring points if the two orientations have a specified direction (e.g. $\langle 0001 \rangle$) in coincidence within a specified tolerance.

Axis/angle – A line segment is drawn between two neighboring points if the misorientation between the two points is within a given tolerance of a specified ideal misorientation where defined by an axis/angle pair (e.g. 60 degrees about $\langle 111 \rangle$).

Grain – A line segment is drawn between two neighboring points if they belong to two different grains.

Phase – A line segment is drawn between two neighboring points if they belong to two different phases.

CSL – A line segment is drawn between two neighboring points if they are within a given tolerance of specified CSL (coincident site lattice boundary).

Shape ellipses – An ellipse is fit to each grain in the map and then overlaid on the map.

Reconstructed – A line is fit to a set of "grain" boundary line segments and drawn on the map.

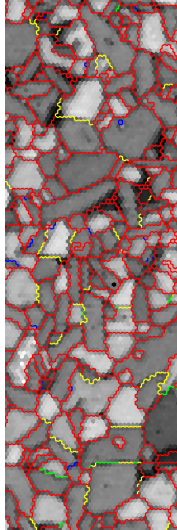
Reconstructed twins – Reconstructed boundaries can be colored according to their fit to specific twin boundary criteria – both in terms of misorientation and twin plane alignment.

Plane traces – A set of lines corresponding to the traces of specified planes are drawn in each grain.

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Boundaries: rotation angle



Rotation angle – This is the most fundamental boundary type. A line segment is drawn between two neighboring points if the misorientation falls in a given range (e.g. between 5 and 15 degrees)

Gray Scale Map Type: Image Quality
11.7...127.2 (11.7...127.2)

Color Coded Map Type: <none>

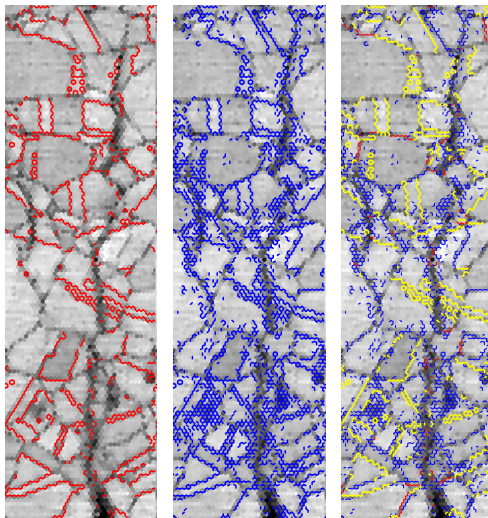
Boundaries: Rotation Angle					
	Min	Max	Fraction	Number	Length
—	3°	6°	0.015	64	18.48 microns
—	6°	9°	0.014	60	17.32 microns
—	9°	15°	0.072	311	89.78 microns
—	15°	180°	0.887	3807	1.10 mm

*For statistics - any point pair with misorientation exceeding 2° is considered a boundary
(total number = 4293, total length = 1.24 mm)

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Boundaries: axis/angle



Rotation angle – A line segment is drawn between two neighboring points if the misorientation falls in a given range. ($60^\circ \pm 5^\circ$ in this example.)

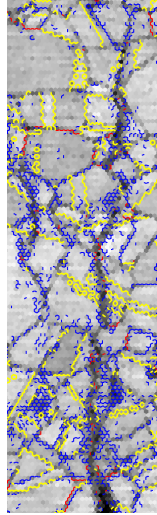
Rotation axis – A line segment is drawn between two neighboring points if the two orientations have specified directions aligned within a specified tolerance. ($\langle 111 \rangle$ crystal directions aligned within 5° of each other in this example.)

Axis/angle – A line segment is drawn between two neighboring points if the misorientation between the two points is within a given tolerance of a specified ideal misorientation where defined by an axis/angle pair. (within 10° of 60° about $\langle 111 \rangle$ in this example.)

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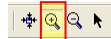
Practice session



- ➔ Open SCCcrack.osc
- ➔ Create an IQ map using QuickGen button



- ➔ Open Map Properties (right mouse click on map)
- ➔ Add rotation angle boundaries
 - Range: 55-65 degrees
 - Color: red (segment tab)
- ➔ Add rotation axis boundaries
 - Axis [uvw]: 111
 - Tolerance: 5 degrees
 - Color: blue
- ➔ Add axis-angle boundaries
 - Tolerance: 10 degrees
 - Angle: 60 degrees
 - Axis [uvw]: 111
 - Color: yellow
- ➔ Extra: change the widths to 9, 5 and 1
 - Zoom in to see effect



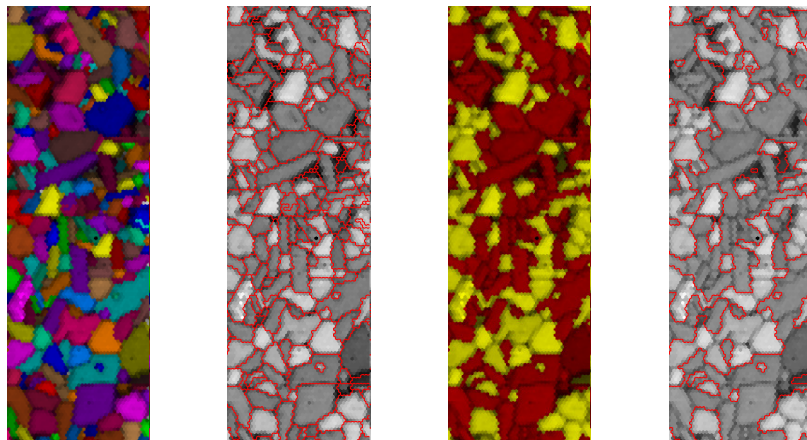
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Boundaries: grains & phase

Grain: A line segment is drawn between two neighboring points if they belong to two different grains.

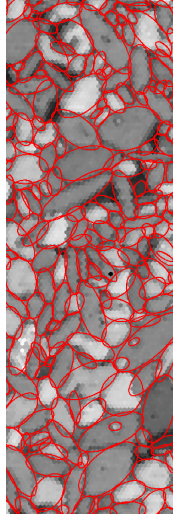
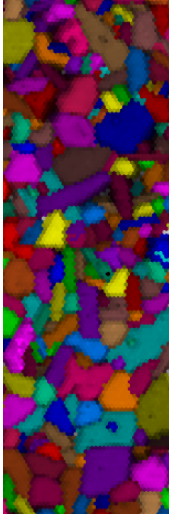
Phase: A line segment is drawn between two neighboring points if they belong to two different phases.



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Boundaries: grain ellipses

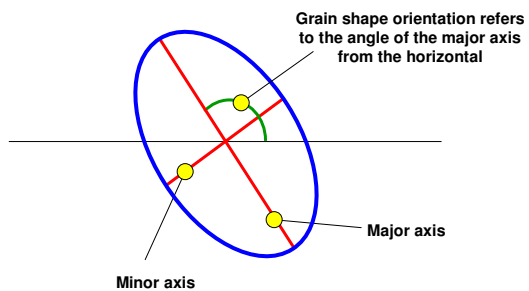
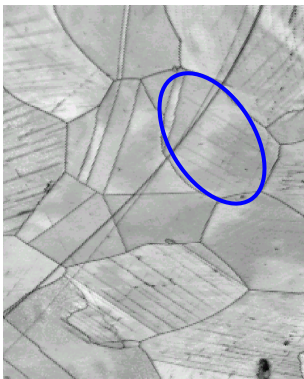


Shape ellipses – An ellipse is fit to each grain in the map and then overlaid on the map. This works best for grains that are not touching the edges of the scan area.

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Boundaries: grain ellipses



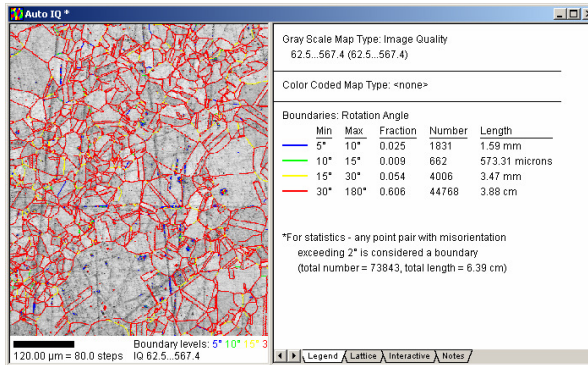
Grain shape aspect ratio is the length of the minor axis divided by the length of the major axis

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Boundaries: minimum misorientation

The minimum boundary misorientation is part of the Settings>Preferences dialog. It is needed for calculations of boundary fractions. For example, if the fraction of boundaries between 5 and 15 degrees misorientation is desired, this fraction is calculated by dividing the number of boundary segments with misorientations between 5 and 15 degrees by the number of boundary segments with misorientations greater than the “minimum boundary misorientation”. Thus, the choice will affect the fractions calculated.



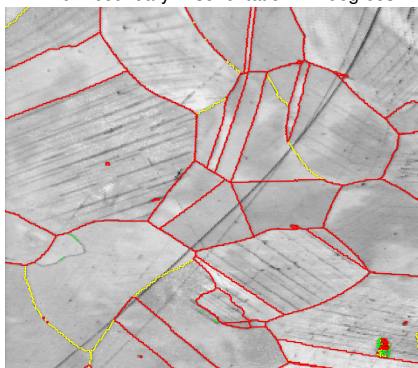
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EDAX TSC

Boundaries: minimum misorientation

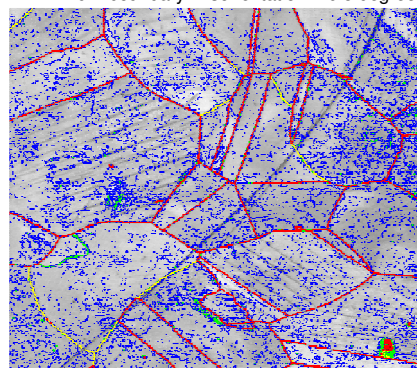
In addition, only boundary segments with misorientations exceeding the “minimum boundary misorientation” will be drawn.

minimum boundary misorientation = 2 degrees



100.00 µm = 100.0 steps Boundary levels: 0.5° 1° 5°
IQ 18.4...207.5

minimum boundary misorientation = 0.5 degrees



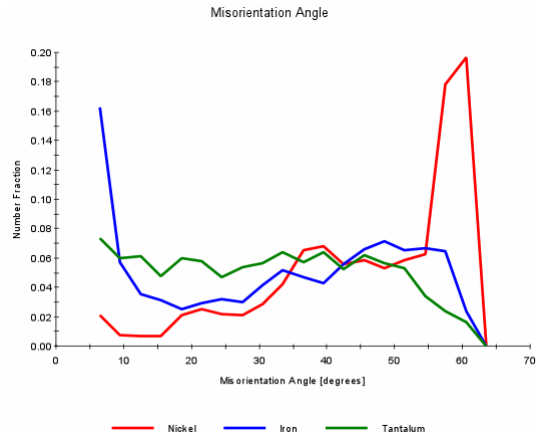
100.00 µm = 100.0 steps Boundary levels: 0.5° 1° 5°
IQ 18.4...207.5

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EDAX TSC

Charts

Image Quality
 Confidence Index
 Video Signal
 Grain Size (points)
 Grain Size (diameter)
 Grain Size (area)
 Grain Shape Orientation
 Grain Shape Aspect Ratio
 Grain Average IQ
 Grain Average CI
 Grain Average Video Signal
 Grain Orientation Spread
 Grain Average Misorientation
 Kernel Average Misorientation
 Crystal Orientation
 Crystal Direction
 Taylor Factor
 Schmid Factor
 EDS
 Phase
 Misorientation Angle
 CSL Boundaries
 Boundary Normal Angle
 Pole Plot
 Pseudo Rocking Curve
 Misorientation Profile
 Texture Fiber
 Texture Index



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EDAX TSE

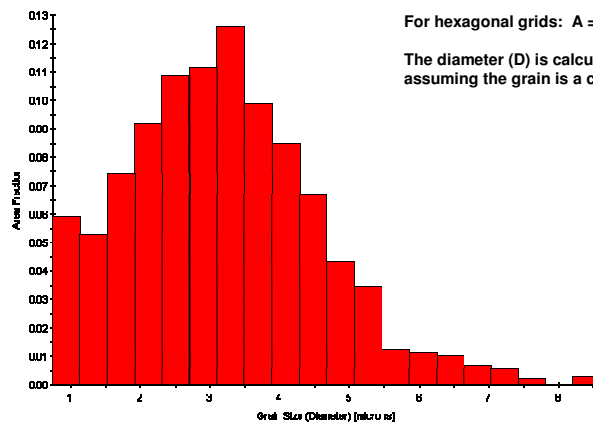
OIM: grain size

The area (A) of a grain is the number (N) of points in the grain multiplied by a factor of the step size (s).

For square grids: $A = Ns^2$

For hexagonal grids: $A = N\sqrt{3}/2s^2$

The diameter (D) is calculated from the area (A) assuming the grain is a circle: $D = (4A/\pi)^{1/2}$.



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EDAX TSE

Charts: grain size

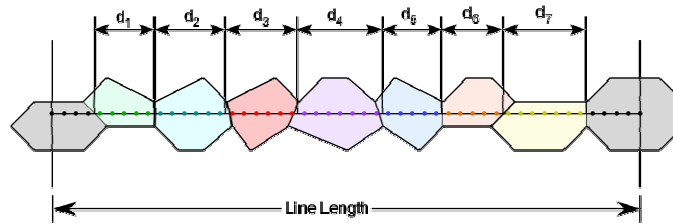
Grain Size (points): Number of measurement points in a grain

Grain Size (diameter): The diameter (D) is calculated from the area (A) assuming the grain is a circle: $D = (4A/\pi)^{1/2}$.

Grain Size (area): The area (A) of a grain is the number (N) of points in the grain multiplied by a factor of the step size (s).
For square grids: $A = Ns^2$. For hexagonal grids: $A = N\sqrt{3}/2s^2$

Grain Size (ASTM): $G = -6.64 \cdot \log_{10}(D) - 2.95$ where D is the grain diameter given in mm

Grain Size (Intercept): Vertical or linear intercepts



Number of grains = 7

Number of intercepts = 8

$$\text{Average diameter} = \bar{d} = \frac{1}{7} \sum_{i=1}^7 d_i$$

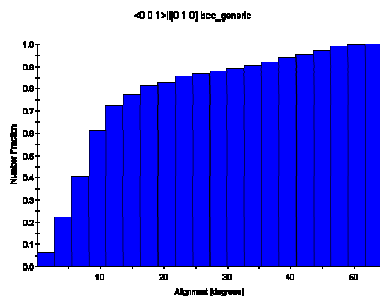
$$\text{Grains per mm} = \bar{d}^{-1}$$

35

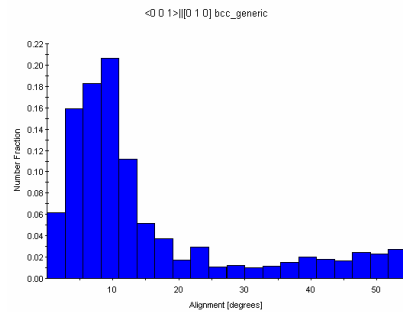
EDAX TSE

Charts: crystal direction

Cumulative



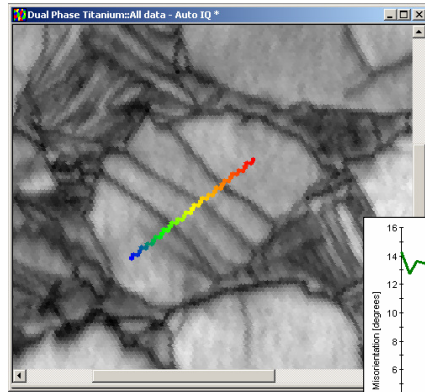
Non-Cumulative



36

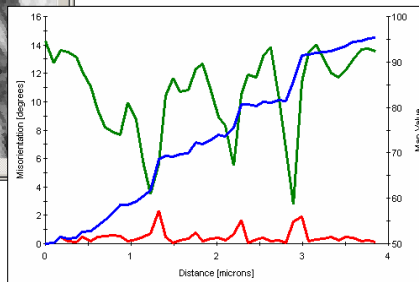
EDAX TSE

Charts: misorientation profile



❖ In addition to the standard misorientation functions, show the map values as well.

- Image Quality
- Point-to-point misorientation
- Point-to-origin misorientation

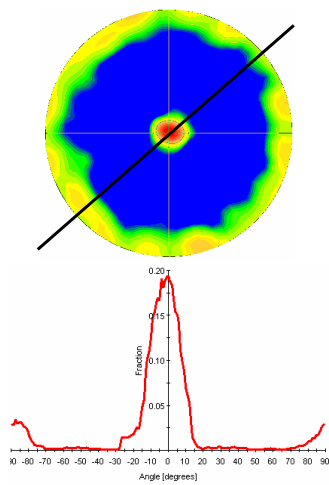


37

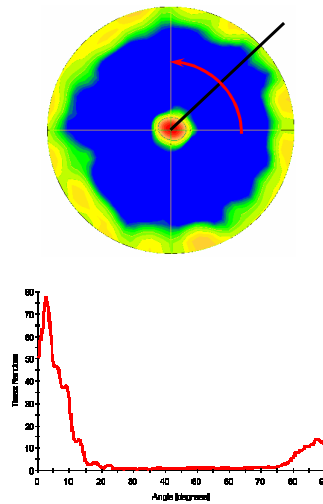
EDAX TSE

Charts: pole plot & pseudo rocking curve

Pseudo Rocking Curve



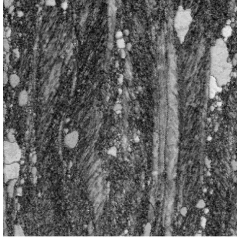
Pole Plot



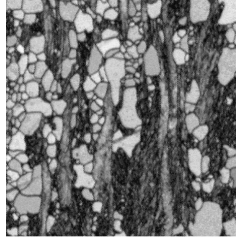
38

EDAX TSE

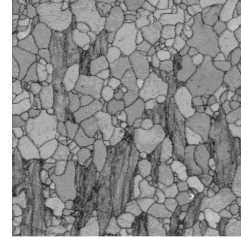
Charts: boundary density



37.50 μm = 50.0 steps IQ 12.5..54.8



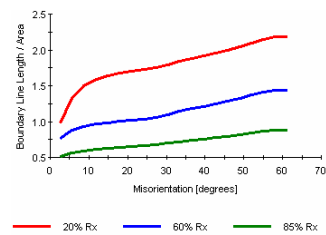
37.50 μm = 50.0 steps IQ 10.4..57.8



37.50 μm = 50.0 steps IQ 11.5..67

In OIM we have considerably more information on the true character of boundaries, it is a shame to discard such information in comparing structures. We have implemented a new chart called a "Boundary Density" chart. The Boundary Density chart shows the line length of boundaries divided by the area as a function of misorientation. It provides a means for linking the boundary structure and grain size together without having as many parameters involved (i.e. boundary tolerance angle, min grain size...).

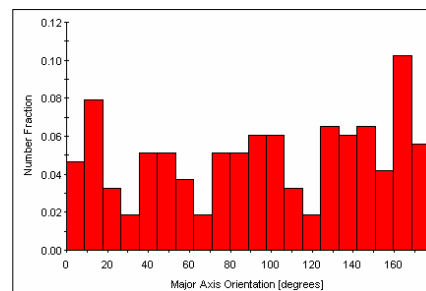
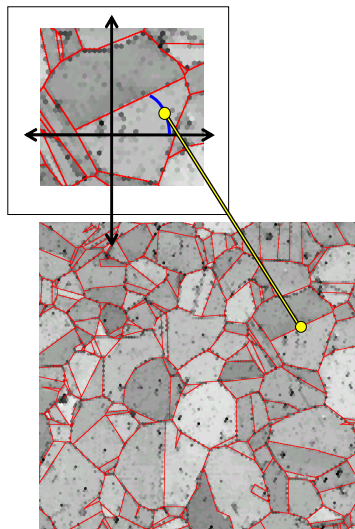
Partially recrystallized low carbon steel.



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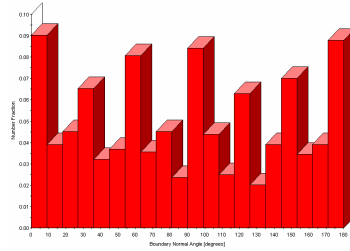
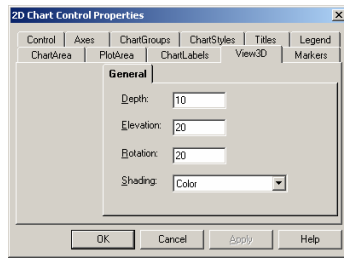
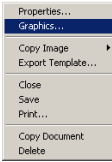
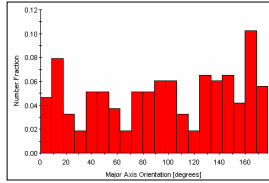
Charts: boundary angle



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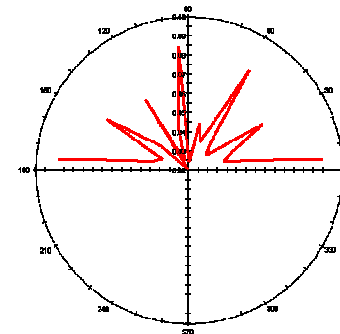
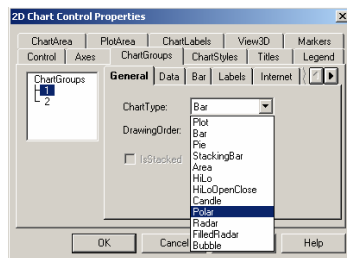
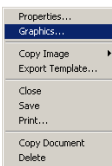
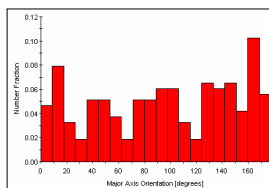
Charts: graphics



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Charts: graphics

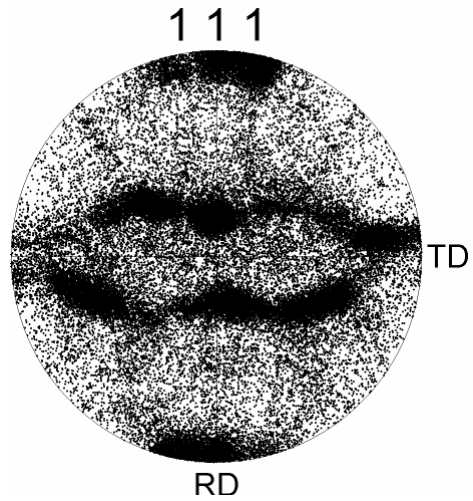


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EDAX TSE

Discrete plots

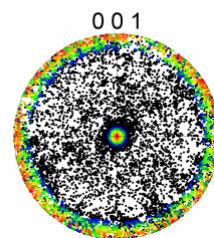
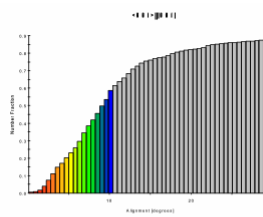
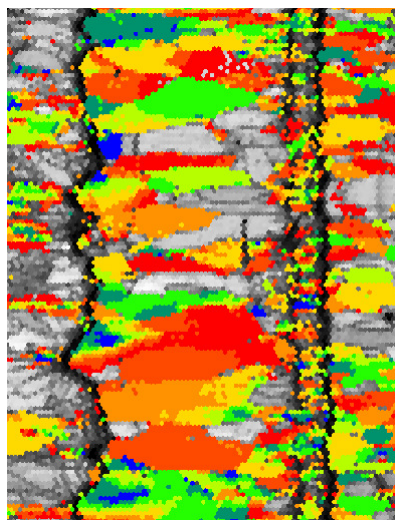
Pole Figure
Inverse Pole Figure
Euler Space
Bunge
Kocks
Roe
Rodrigues Orientations
Rodrigues Misorientations
Axis/Angle Misorientations



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EDAX TSE

Highlighting

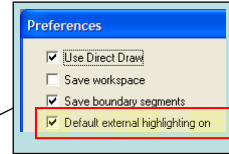


44

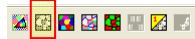
EDAX TSE

Practice session - highlighting

- ➔ Open CVDTa.osc
- ➔ Go to the Settings menu and select Preferences and make sure highlighting is defaulted on.
- ➔ Create an IQ map using the QuickGen button



- ➔ Create a (001) pole figure using the QuickGen button



- ➔ Create a Crystal Direction Chart
 - Set the Sample Direction to RD=0, TD=1, ND=0
- ➔ Next Slide...

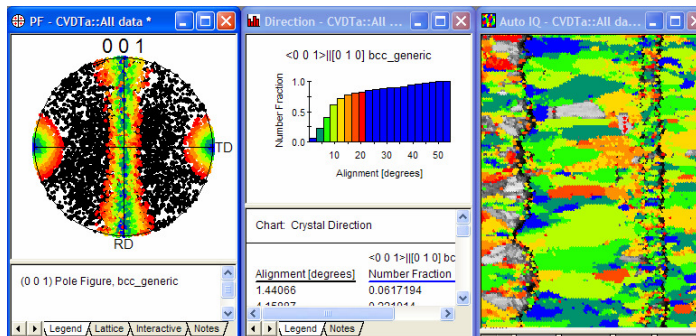
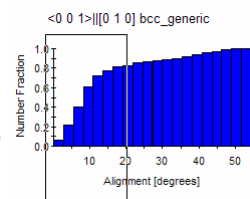


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EDAX TSE

Practice session – highlighting continued

- ➔ Drag a box on the chart
 - Make sure you start inside chart area
- ➔ Note Highlighting Effects
 - Change the pole figure markers to a larger size to see effects better (right mouse click to select properties, press “Edit >>” and select point tab.



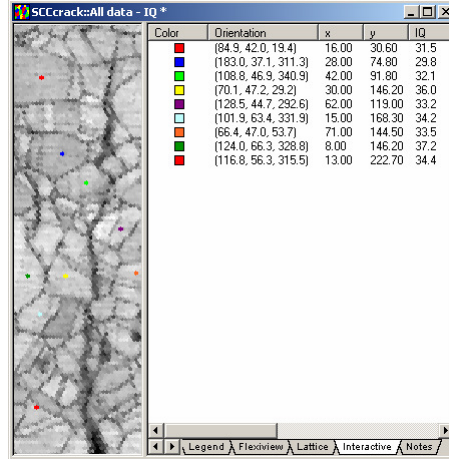
46

EDAX TSE

Highlighting



Record: record highlighting results in the Interactive pane of the map window.



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EDAXTM TSE

Highlighting: toolbar



Undo previous highlight.
Redo deleted highlight.
Clear all highlights.

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EDAXTM TSE

Practice session - highlighting

- Open SCCcrack.osc
- Create an IQ map
- Turn on “record” button for highlighting



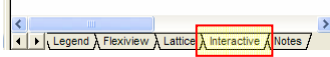
- Select the single point highlighting mode



- Click on a few points in the map
- Switch colors in-between clicks



- Switch to the “interactive” pane of the map window



- Note the items recorded in the table
 - Right mouse click to show preferences for interactive table
- Follow along with subsequent highlighting examples

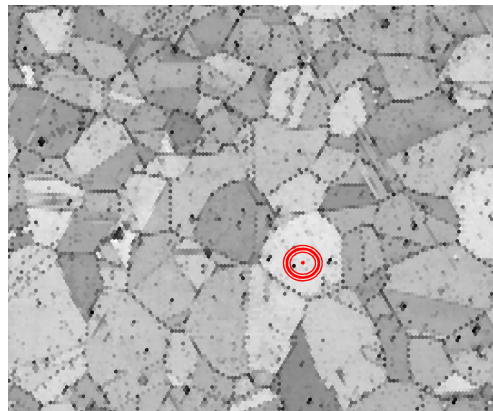
49

EDAX[™] TSE

Highlighting: toolbar



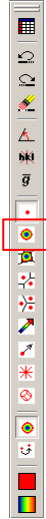
Highlight the point clicked.



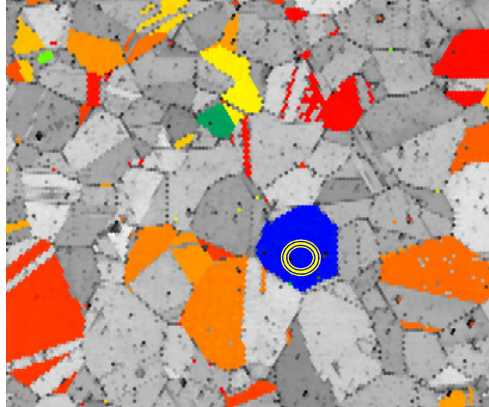
50

EDAX[™] TSE

Highlighting: toolbar



Highlight all orientations within a given tolerance of the point clicked.



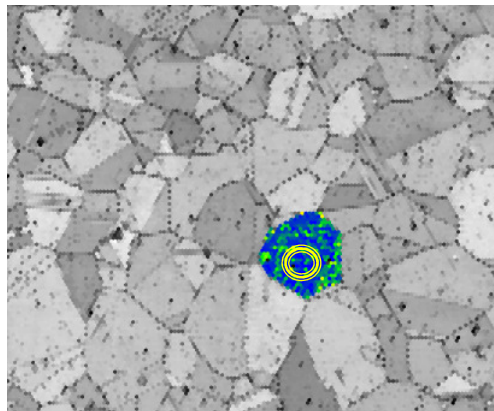
51

EDAX[™] TSE

Highlighting: toolbar



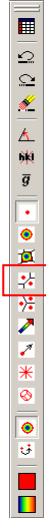
Highlight all orientations within grain and color them according to the angular distance from the point clicked.



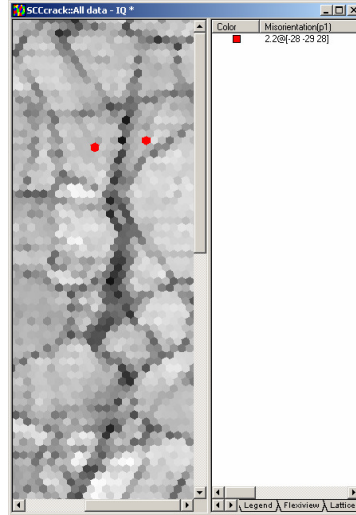
52

EDAX[™] TSE

Highlighting: toolbar



Calculate misorientation
between two points.



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EDAX TSC

Highlighting: toolbar



Calculate misorientations
at a triple junction

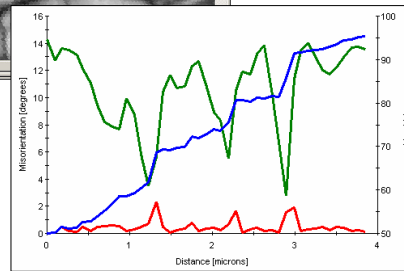
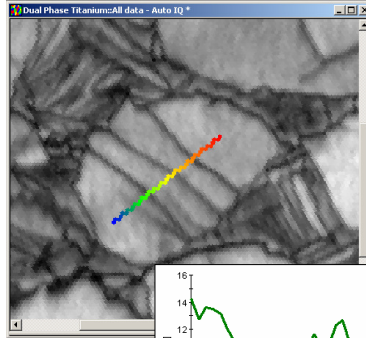
54

EDAX TSC

Highlighting: toolbar



Misorientation profile:
Creates a chart plotting
misorientations along a
line.



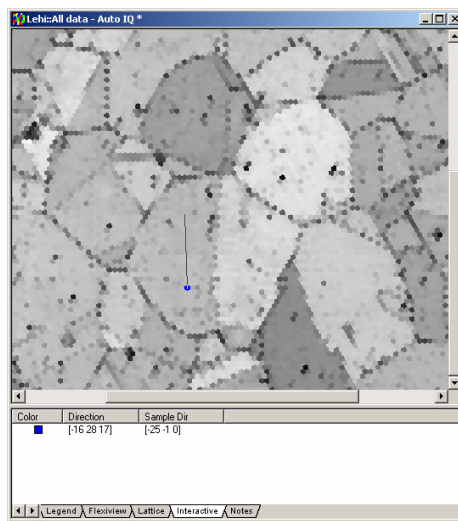
EDAX TSE

55

Highlighting: toolbar



Displays a crystal
direction parallel to
a line.



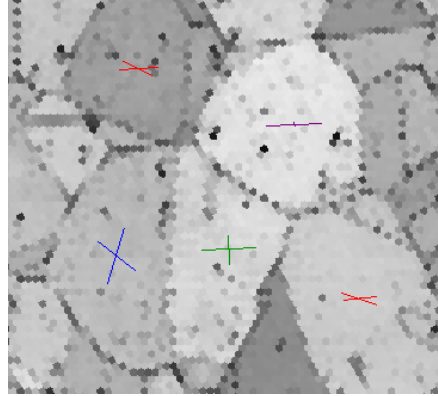
EDAX TSE

56

Highlighting: toolbar



Draw traces of a specified plane at a point.



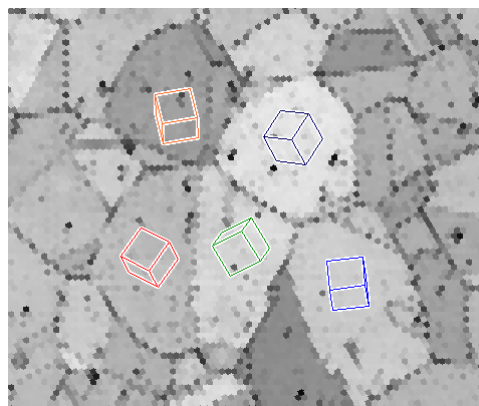
57

EDAXTM TSC

Highlighting: toolbar



Draw a unit cell representing the orientation at a point.



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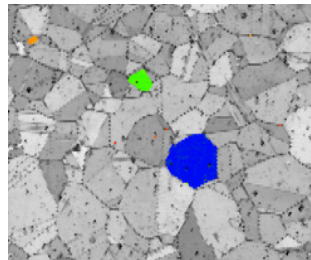
EDAXTM TSC

Highlighting: toolbar

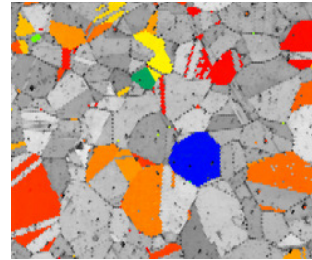


Angular tolerance used for various highlighting modes.

Tolerance = 15°



Tolerance = 30°



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EDAX TSC

Highlighting: toolbar



Define the (hkl) for the plane traces.

Trace Plane Indices

Define the plane

h	k	l
1	1	-2

Normal vs. Trace

Show the plane trace

Show the plane normal

The longer the trace/normal the closer the plane is to perpendicular to the sample surface.

OK Cancel Apply Help

60

EDAX TSC

Highlighting: toolbar



Use the average grain orientations when calculating misorientations.

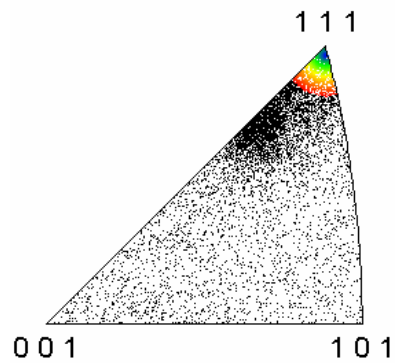
61

EDAXTM TSL

Highlighting: toolbar



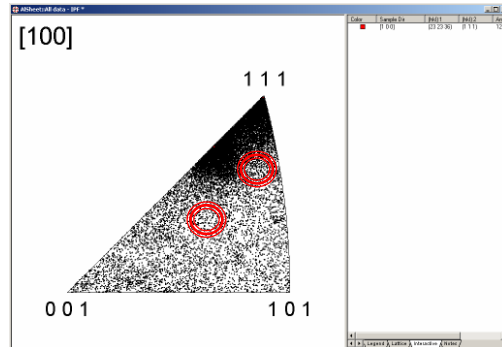
Find all points in the scan within the angular tolerance of the point clicked in discrete plot.



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EDAXTM TSL

Highlighting: toolbar



Calculate the misorientation between two points in a discrete plot

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EDAX TSC

Highlighting: toolbar

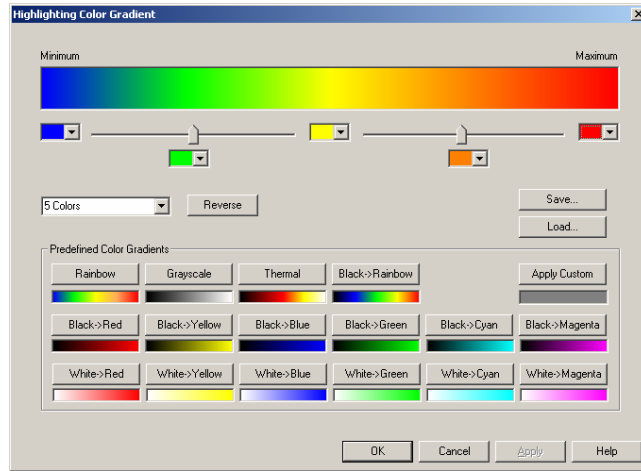


Select a color for individual highlights.

64

EDAX TSC

Highlighting: toolbar

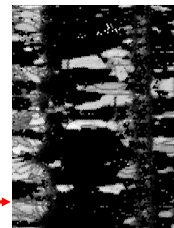
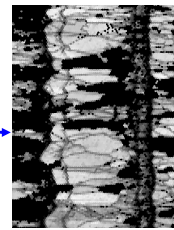
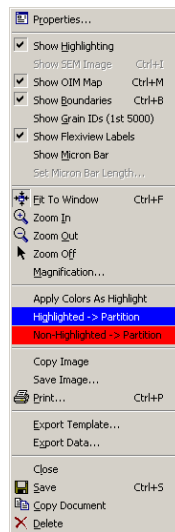
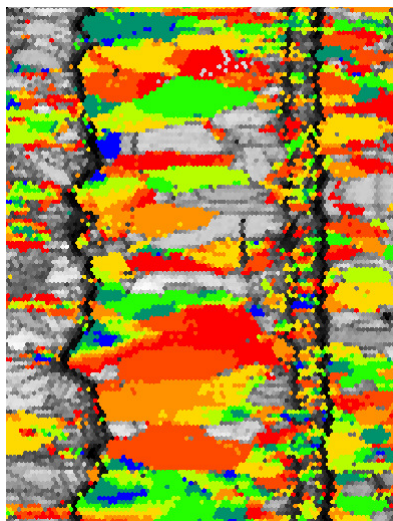


Select a color gradient
multiple point highlights.

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EDAX TSE

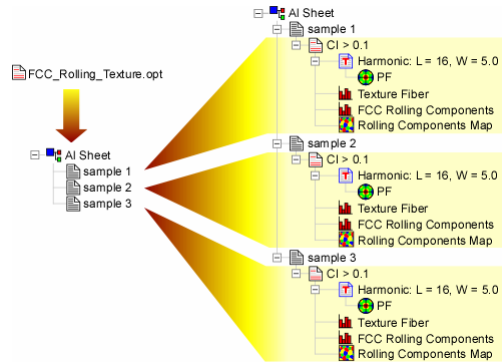
Highlighted → partition



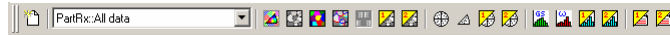
66

EDAX TSE

Templates



User defined templates can be assigned to buttons on the QuickGen Toolbar.

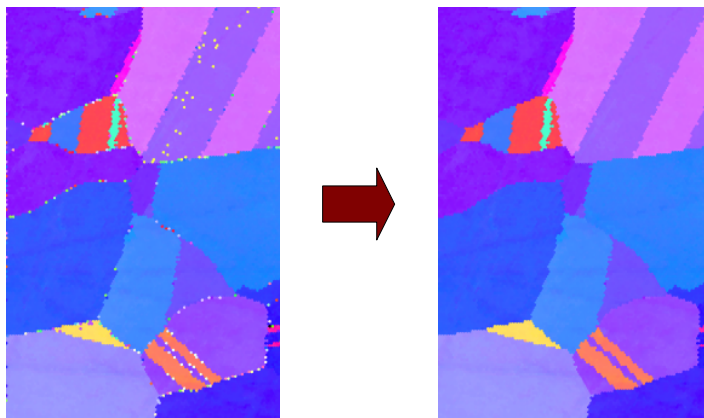


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EDAX[®] TSE

Clean up

Sometimes there are isolated points that are not indexed correctly or at all due to dust particles or pits on the surface. Various clean up schemes are available to assign these points an orientation based on their neighbors.

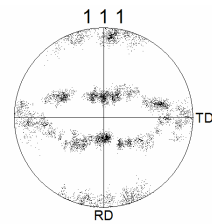
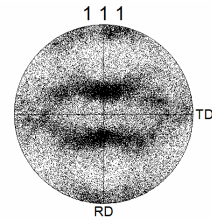
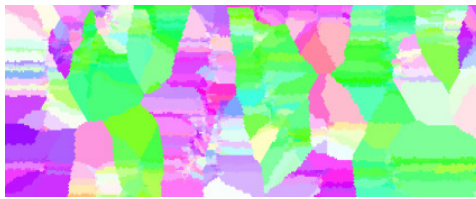
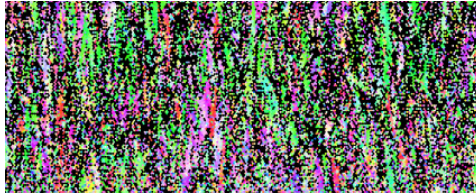


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EDAX[®] TSE

Clean up

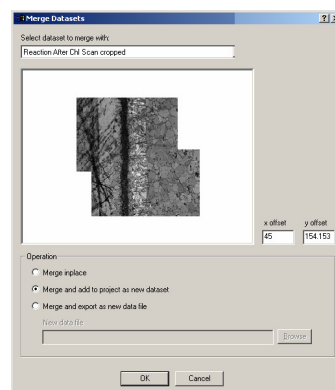
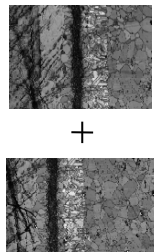
However, it is important to make sure that the resulting microstructure makes sense also watch the **texture** and grain boundary changes.



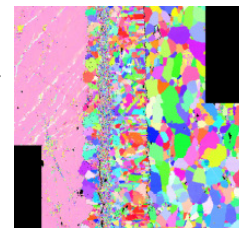
69

EDAXTM TSL

Merge



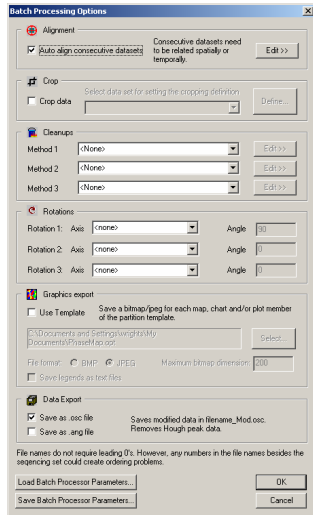
- ❖ Datasets can be merged together.
- ❖ The relative positioning of the datasets is controlled by the user.



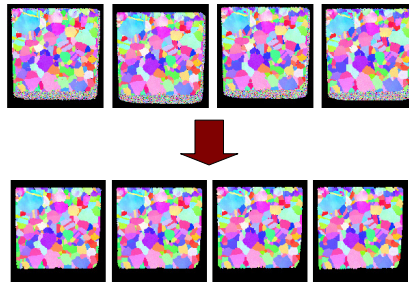
70

EDAXTM TSL

Batch processor



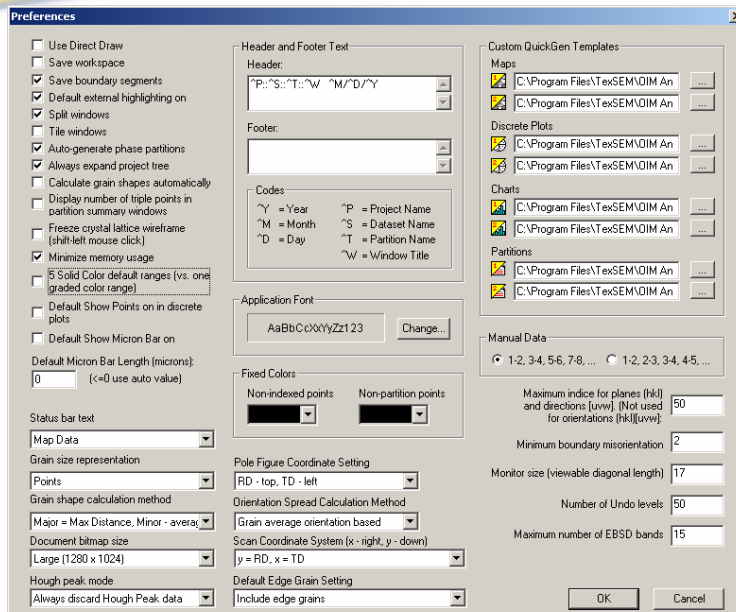
- ❖ A tool to apply a partition template to multiple sets of data.
- ❖ Includes Rotate, Cleanup, Crop, and Export functions.
- ❖ An alignment function to align consecutive sets of data for a sequence of scans, e.g. from serial sectioning or slices in time.



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Preferences



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