Introducing the 2020 Powder Diffraction File™ -

# Diffraction Data You Can Trust

ICDD databases are the only crystallographic databases in the world with quality marks and quality review processes that are ISO certified.

# PDF-4+/Web **Identification** and Quantitation 426,000+ Entries 323,000+ Atomic coordinates **Combines Digitized Powder and Patterns** Single Crystal Data Analyze Molecular Neutron, Electron, X-ray, Synchrotron **Graphics** Data

Standardized Data

More Coverage

All Data Sets Evaluated For Quality

Reviewed, Edited and Corrected Prior To Publication

Targeted For Material Identification and Characterization







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# Diffraction Data You Can Trust

ICDD databases are the only crystallographic databases in the world with quality marks and quality review processes that are ISO certified.

# PDF-4/Axiom **Focused** Identification and Quantitation 92,000+ Entries 63,000+ Atomic coordinates **Combines Cost Effective Powder and** 3 Year Single Crystal License Data Requires **Two Additional** Vendor Seats - Low Cost Software

Standardized Data

All Data Sets Evaluated For Quality

Reviewed, Edited and Corrected Prior To Publication

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# LET OUR TEAM OF EXPERTS HELP YOU TAKE YOUR SKILLS TO THE NEXT LEVEL!



#### Rietveld Refinement & Indexing Clinic:

#### 23 - 27 September 2019

Powder pattern indexing and Rietveld structural refinement techniques are complementary and are often used to completely describe the structure of a material. Successful indexing of a powder pattern is considered strong evidence for phase purity. Indexing is considered a prelude to determining the crystal structure, and permits phase identification by lattice matching techniques. This clinic introduces the theory and formalisms of various indexing methods and structural refinement techniques along with quantitative analysis. One unique aspect of this clinic is the extensive use of computer laboratory problem solving and exercises that teach method development in a hands-on environment.



#### **Practical X-ray Fluorescence Clinic:**

#### 27 April - 1 May 2020

From theory to hands-on exercises, this course offers techniques and skills to improve lab performance. Discover the latest in cutting-edge instruments such as TXRF, hand-held devices, energy dispersive and wavelength dispersive spectrometers through live demonstrations.

The XRF course covers the basics of X-ray spectra; instrumentation design; methods of qualitative and quantitative analysis; specimen preparation and applications for both wavelength and energy dispersive spectrometry. The course emphasizes quantitative methods, use of automated X-ray spectrometers, review of mathematical matrix correction procedures, and new developments in XRF.



#### Fundamentals of X-ray Powder Diffraction Clinic:

#### 1 - 5 June 2020

For the novice with some XRD knowledge or for the experienced with an interest in the theory behind XRD, this clinic offers a strong base for increased lab performance.

The clinic covers instrumentation, specimen preparation, data acquisition and qualitative phase analysis through live demonstrations. It also covers hands-on use of personal computers for demonstration of the latest software including data mining with the Powder Diffraction File (PDF) and use of the powder diffractometer: optical arrangement, factors affecting instrumentation profile width, choice and function of divergence slit, calibration and alignment, detectors, and X-ray optics.



#### Advanced Methods in X-ray Powder Diffraction Clinic:

#### 8 - 12 June 2020

For the experienced XRD scientist, this session offers enhanced analysis skills through intense problem solving, as well as an introduction to the Rietveld Method. The course emphasizes computer-based methods of data collection and interpretation, both for qualitative and quantitative phase analysis.

The advanced clinic covers factors affecting d-spacings of crystals, as well as factors affecting diffraction-line intensities; structure-sensitive properties (atomic scattering and structure factors), polarization effects, and multiplicity. Additionally, the clinic covers specimen-sensitive effects (orientation, particle size), measurement-sensitive effects (use of peak heights and peak areas), and choice of scanning conditions will also be addressed.

Please visit the ICDD website for more information.

#### Register today at WWW.ICDD.COM

**Please note:** A minimum of 10 registrants per course is required, otherwise the course will be cancelled and your registration fee will be refunded. You will be notified of a course cancellation no later than two weeks prior to the start of the course.







#### For More Information Contact:

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#### Location

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# Better TOGETHER

## **SAVE 15% on our New Bundles**

PDF-4+ and JADE Pro 2020 PDF-4/Organics and JADE Pro 2020



## **Phase Identification & Quantitation**

PDF-4+ features more data, higher quality, more content standardized data, and editorial evaluated data reviewed, edited and corrected prior to publication



# Solve Difficult Problems, Get Better Results

PDF-4/Organics database is a highly targeted collection, with special focus on materials used in commercial and regulatory fields. It provides the best of both worlds by including single crystal and powder diffraction data together in a single, edited, and standardized database.



## Scientists Trust Materials Data

JADE Pro Is a powerful, all-purpose powder XRD pattern viewer, processing and analysis program with emphasis on quantification and phase ID.

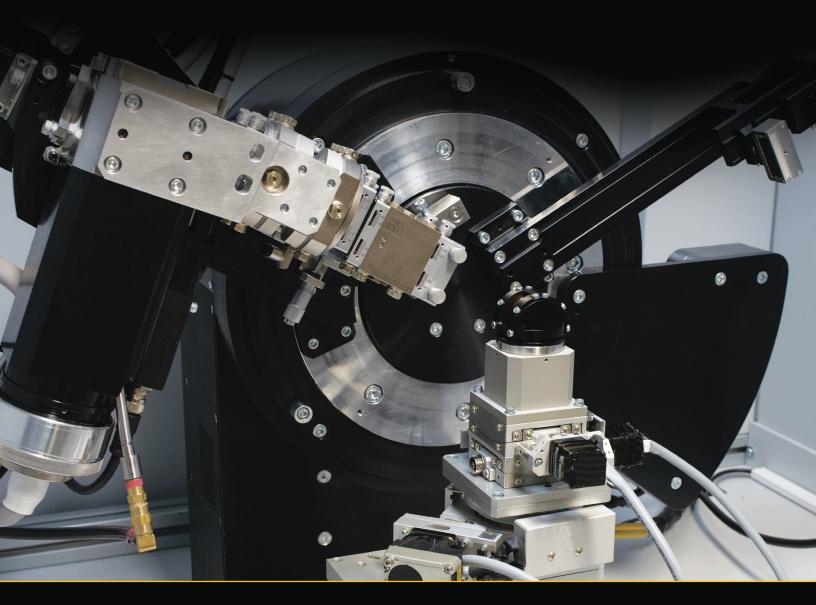




# HIGH-RESOLUTION DIFFRACTION

Customizable high-resolution diffractometers for characterizing thin films and single crystal materials.

- · X-RAY REFLECTIVITY (XRR) · ROCKING CURVES
- · RECIPROCAL SPACE MAPPING



PROTO

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