## 2025-2026 **SALES** CATALOG

The Powder Diffraction File™ (PDF®) & JADE Software

PDF# 00-075-0477 Na<sub>3</sub>V<sub>1.95</sub> Nb<sub>0.05</sub>(PO<sub>4</sub>)<sub>2</sub>O<sub>2</sub>F Sodium Vanadium Niobium Oxide Fluoride Phosphate Wang, J., Yuan, Y., Rao, X., Yang, M., Wang, D., Zhang, A., Chen, Y., Li, Z., Zhao, H.

> ed Profile Raw Diffraction Data

Raman Spectra •

aw diffraction data

Fixed Slit Intensity

d (Å)	I (%)	h	k	1	*	
5.46363	33	1	0	1		
5.31790	73	0	0	2		
4.50293	8	1	1	0		
3,43646	6	1	1	2		

6000 Intensity 2000 2000 0 20 30

Status:

**Environment:** 

Primary

Quality Mark: Star

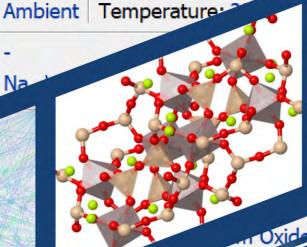


Phase:

ital

Chemical Formula: Na

**Empirical** 









## Our collaborative partners include:

Cambridge Crystallographic Data Centre (CCDC) Cambridge, UK - Cambridge Structural Database (CSD)

## FIZ Karlsruhe - Leibniz Institute for Information Infrastructure (FIZ)

Karlsruhe, Germany - Inorganic Crystal Structure Database (ICSD)

## National Institute of Standards and Technology (NIST)

Gaithersburg, Maryland, USA

Material Phases Data System (MPDS)

Rome, Italy - Linus Pauling File (LPF)

## PDF Databases - Why You Need The Best!

- More data, higher quality, more content, many types of solid-state reference data
- The world's largest solid-state collection of minerals, metals and alloys, polymers, active pharmaceuticals, and commercial materials
- The world's largest and most diverse (by formula) collection of inorganic structures
- The world's largest collection of modulated structures
- Extensive bibliographic collections thousands of journals, hundreds of thousands of authors
- Standardized data, nomenclature, and structural classifications plus brand names, trade names, mineral names, zeolite names, and common names of commercial materials
- Identify materials accurately by utilizing the PDF's extensive subfile system that classifies materials by chemistry and application
- New data is contributed annually by highly qualified grantees from around the world
- PDF products are compatible with major diffraction vendor's software



Quality Management System Certified to ISO 9001:2015 by DEKRA

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#### THE INTERNATIONAL CENTRE FOR DIFFRACTION DATA (ICDD®)

is a non-profit scientific organization dedicated to collecting, editing, publishing, and distributing powder diffraction data for the identification of materials. The membership of the ICDD consists of worldwide representation from academe, government, and industry.

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Grant-in-Aid	

#### ONLINE RESOURCES

Visit www.icdd.com. Learn from reference articles, technical bulletins, tutorials, and webinars. Some tutorials and webinars focus on the capabilities of the PDF database, while others illustrate methods used to analyze pharmaceuticals, polymers, and minerals. It is free to search and browse documents from ICDD's Pharmaceutical Powder X-ray Diffraction Symposium (PPXRD) and the ICDD Spring Meetings. You get open access to over 1,000 full publications from Advances in X-ray Analysis (AXA) - the proceedings of the Denver X-ray Conference® (DXC). Try the Powder Diffraction File (PDF) Search and search using chemical name, formula, and elements to suggest which ICDD database product is right for you.



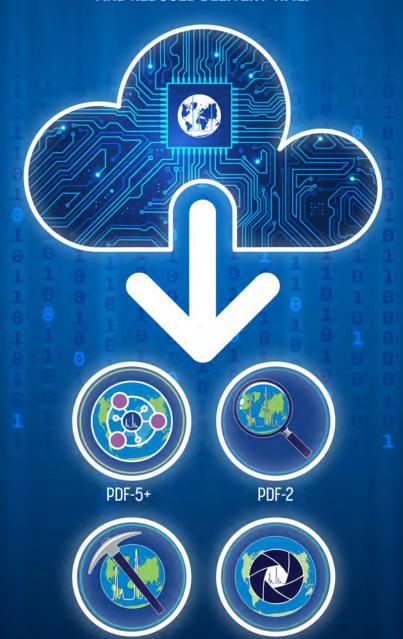


# THE POWDER DIFFRACTON FILE IS AVAILABLE FOR

# **DOWNLOAD!**

GET YOUR 2026 PDF DATABASE BY DOWNLOADING FROM THE CLOUD

BENEFIT FROM NO SHIPPING COST AND REDUCED DELIVERY TIME!





## LICENSE OPTIONS TO FIT YOUR BUDGET!

- Extended multi-year license terms -Let us know what you need!
- Customized license for a 1 to 9-year license term
- 15% discount for new PDF-5+ and IADE Pro Bundles
- One-stop shopping: align PDF and JADE renewals for one billing and shipment
- PDF-2 10-year license prepaid package
- PDF-5+Scholar -10-year license for Academics
- PDF-4/Axiom 6-year and 9-year license - prepaid packages
- SAVE time and money with ICDD's PDF-5+ and PDF-4/Minerals Subscriptions
  - ICDD's subscriptions keep your database up-to-date with renewal pricing
  - ICDD offers automatic PDF-5+ and PDF-4/Minerals renewal shipments/download renewals for convenience
  - Can't accept automatic shipments?
     ICDD will send a PDF-5+ or
     PDF-4/Minerals renewal reminder to keep you up-to-date with your subscription

PDF-4/AXIOM

PDF-4/MINERALS

## **WHAT'S NEW FOR 2025-2026**



**PDF-5+** *The* Powder Diffraction Standard



PDF-5+ Server Edition
Extend the Power of Your Database –
Built for Teams of Five or More



PDF-4/Axiom
Quality plus Value



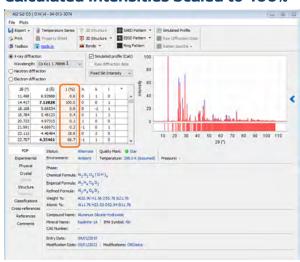
PDF-4/Minerals
Comprehensive Mineral Collection



**PDF-2**Phase Identification + Value

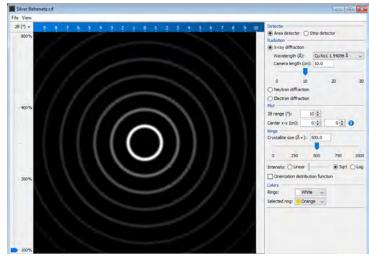
### Introducing Release 2026 of the Powder Diffraction File

#### **Calculated Intensities Scaled to 100%**

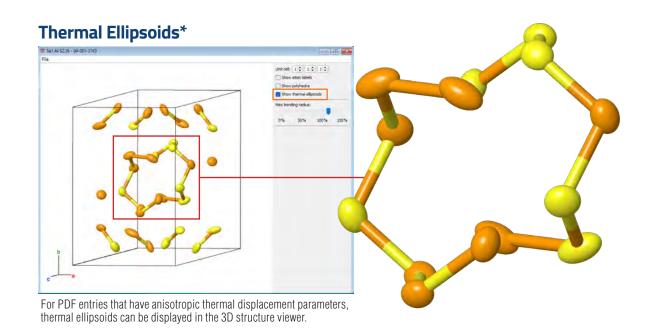


- For historical reasons, intensities for calculated diffraction patterns were scaled to 1000 using integers
- Now intensities are scaled to 100% using floating-point numbers

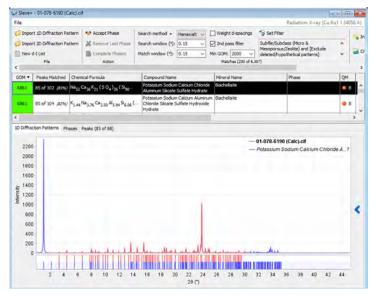
#### Simulate 2D Ring Patterns from CIFs\*



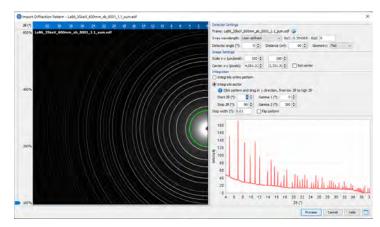
In addition to simulating 2D ring patterns from PDF entries, 2D ring patterns can now be simulated from a user imported Crystal Structure \*.cif file.



#### **New Sleve+ Features**



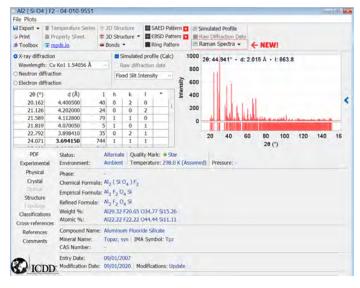
- Combine the search history with common filters
- · Deep Search option
  - When selected, the current search will use all the permutations of 8 strongest search peaks to 3 strongest database peaks (instead of just the strongest database peak)
- New X-axis range option
  - Full Range
  - Restrict Range (to experimental data)
- Hide phase plots using checkboxes\*



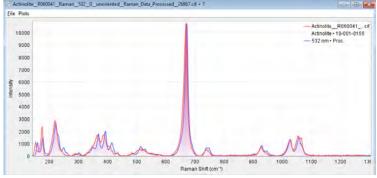
- Import 2D diffraction patterns from ESRF files (\*.edf)\*
- Simplified sector integration with 2D diffraction patterns\*

#### **Raman Database**

With the increasing interest in identifying minerals using Raman spectroscopy, it is time for a unified effort to publish a curated Raman database. The International Centre for Diffraction Data (ICDD), with over 8 decades of experience with reference database science and an extensive history with the minerals community, is uniquely suited to this task. Raman data is available in PDF-5+ 2026, PDF-4/Minerals 2026, and PDF-4/Axiom 2026.



#### **New Raman Features**



- Raman Analysis Graph\*
  - The user can import Raman spectra from a file and add Raman spectra from the PDF database into a graph for visual analysis
- · Raman-Diffraction Search\*
  - An option has been added to the Raman search that will only include Raman entries that are cross-referenced to X-ray diffraction data
- New Raman Display Fields\*
  - · Literature reference
  - Database comments

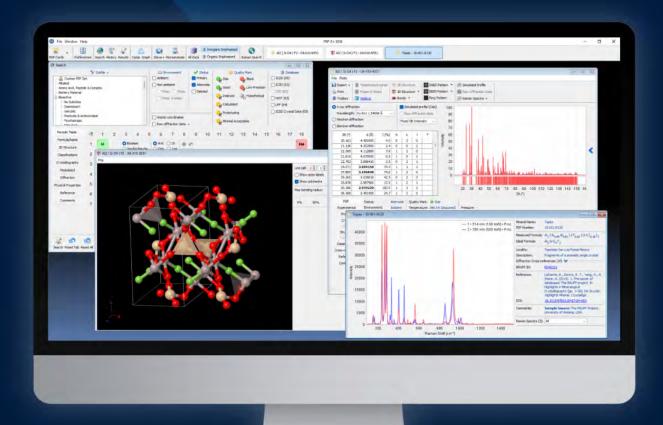
## PDF-5+2026 THE POWDER DIFFRACTION STANDARD

Inorganic & Organic entries are now combined into one powerful database with 1,126,200+ entries.



#### www.icdd.com/pdf-5

- 651,200+ entries with atomic coordinates
- Features 472,300+ entries for inorganic materials and 657,700+ entries for organic materials
- All entries have digital patterns for use in total pattern analysis
- 1,022,600+ entries have I/I<sub>c</sub> values for quantitative analysis by Reference Intensity Ratio
- All entries are stored in a standardized format for easy search and interpretation
- All entries go through a rigorous editorial process to ensure quality

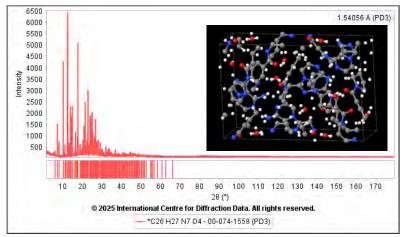




## PDF-5+ 2026

### The Powder Diffraction Standard

PDF-5+ combines the world's largest sources of inorganic and organic diffraction data from crystals and powders into a single database.



Calquence®, acalabutinib dihydrate Form III, used to treat various types of non-Hodgkin lymphoma.

## **(7)** KEY POINTS

- Rapid phase identification designed to support automated quantitative analyses
- Features 1,126,200+ entries with over 651,200+ having atomic coordinates
  - > 472,300+ entries for inorganic materials
  - > 657,700+ entries for organic materials
- Integrated data mining software
- All entries have digital patterns for use in total pattern analysis
- More than 1,022,600+ entries have I/I<sub>c</sub> values for quantitative analysis by Reference Intensity Ratio (RIR) method
- Combines powder and single crystal data
- Includes crystalline, semi-crystalline, and amorphous solid-state materials
- World's largest and most diverse collection of structures for inorganic and organic materials
- Extensive data mining is facilitated with 118 display fields coupled with 85 searches
- LPF entries have a link to mpds.io
- One (1) year license serviced by subscription

**PDF-5+ Purchase Options** 

PDF-5+ 2026	USB CODE	DOWNLOAD CODE	2025-2026 LIST \$	2025-2026 Academic \$
New - Single License	P5S26	P5L26	\$9,395	\$6,265
Renewal from 2025 to 2026	P5RS26	P5RL26	\$2,115	\$1,375
PDF-5+ and JADE Pro bundle*	'		'	
New - Single License	BUJPP5S26	BUJPP5L26	\$16,310	\$11,465
*Purchase a new license for PDF-5+ and JADE Pro				
PDF-5+ 2026 Multi-year License*	'		'	
Renewal – 3 year License	P5S27MR29	P5L27MR29	\$5,710	\$3,715
Renewal – 5 year License	P5S27MR31	P5L27MR31	\$8,990	\$5,845
PDF-5+ Scholar 2026 (10-Year License Package)*				
PDF-5+ New — Single License	P5S26	P5L26		\$6,265
Scholar Multi-year	P5SPS27MR31	P5SPL27MR31	NA	\$5,845
	·	· ·		\$12,110

#### Save Money! Take advantage of ICDD's Site License - Request a quote today!

Save money. Take advantage of 1000 3 Site Election - nequest a quote today.					
PDF-5+ 2026 Site License*	USB CODE	DOWNLOAD CODE	2025-2026 LIST \$	2025-2026 ACADEMIC \$	
New - Master License	P5S26M	P5L26M	\$9,395	\$6,265	
New - additional License, each	P5S26C	P5L26C	\$2,115	\$1,375	
Renewal - Yearly Master License	P5RS26M	P5RL26M	\$2,115	\$1,375	
Renewal - additional License, each	P5RS26C	P5RL26C	\$855	\$590	
PDF-5+ 2026 Site/Multi-year License*					
Renewal - 3 year Master Site License	P5S27MR29M	P5L27MR29M	\$5,710	\$3,715	
Renewal - 3 year additional Site License, each	P5S27MR29C	P5L27MR29C	\$2,565	\$1,770	
Renewal - 5 year Master Site License	P5S27MR31M	P5L27MR31M	\$8,990	\$5,845	
Renewal - 5 year additional License, each	P5S27MR31C	P5L27MR31C	\$4,275	\$2,950	
PDF-2 to PDF-5+ Conversion	·				
PDF-2 2025 to PDF-5+ 2026	P2525RSU26	P2525RLU26	\$2,115	\$1,375	
PDF-2 2024 to PDF-5+ 2026	P2524RSU26	P2524RLU26	\$4,230	\$2,750	
PDF-2 2023 to PDF-5+ 2026	P2523RSU26	P2523RLU26	\$6,345	\$4,125	
PDF-2 2022 to PDF-5+ 2026	P2522RSU26	P2522RLU26	\$8,460	\$5,500	
	•	The second secon			

<sup>\*</sup>Please see product policies on pages 18 and 19



## **Extend the Power of Your Database**Built for Teams of Five or More

Workplaces and work habits have evolved from just a few years ago, and this database licensing model provides the flexibility that large organizations and academics need to meet the varying demands of clients, teams, and students.

Install on your server and as many computers as you wish. A 5-User license provides access to 5 computers at a time and floats among an unlimited number of systems. The Server license can be expanded to 10, 15, or more users. Deployment to new hires is seamless. When it's time for the latest database, your renewal is simplified with one single product. Ensure everyone in your organization is using the latest set of data with one shared floating 5 or more seat license. Additional benefits occur when access is granted to occasional use individuals without any added cost to your bottom line.

Your IT department will be happy as updates are streamlined, all while offering savings for your budget. If you need to provide access to 5 or more systems, please contact us at <a href="mailto:sales@icdd.com">sales@icdd.com</a> to consider how our new Server License would work best for your team.



# PDF-5+ 2026 The Powder Diffraction Standard

PDF-5+ combines the world's largest sources of inorganic and organic diffraction data from crystals and powders into a single database.

PDF-5+ contains 472,300+ inorganic and 657,700+ organic entries to ensure that you have the data you need for all of your phase identification needs. Every entry is editorially reviewed and assigned quality marks, so the curated PDF-5+ database allows users to do quantitative analysis by any of three methods: Rietveld Analysis, Reference Intensity Ratio (RIR) method, or Whole Pattern Fitting.



## **JADE Pro**

## The Analysis Software Professionals Choose

You need data you can trust, AND you need to find the best way possible to analyze your data. Extend the power of your ICDD PDF-5+ Server Edition by including the highly regarded data analysis software JADE Pro. We created JADE to provide independent, unbiased results for peak search, whole pattern fitting, and Rietveld. JADE gets daily improvements and has grown to include an enormous list of valuable tools for your materials research and exploration.

The power of JADE combined with the PDF-5+ Server Edition means a team can easily share a 5-seat license. The Server Edition can result in cost savings for your organization and a chance for you to extend the power to your team by providing access to BOTH the PDF database and JADE to as many individuals as you need. The database and software products offer the same functionality as the individual products. In coupling them in a 5 or more user server license, important resources are available to more team members, and everyone is in sync, all while reducing costs.



- All-inclusive Everything in JADE Standard and so much more
- Phase ID (Search/Match)
- Batch processing Whole Pattern Fitting (WPF) and Rietveld
- One-Click-Analysis<sup>™</sup> for Whole Pattern Fitting (WPF)
- Pattern Indexing (All Crystal Systems)
- Rietveld Structure Refinement (Atomic Parameters)
- ab Initio Tools (Charge Flipping +)
- Cluster Analysis of Powder Patterns
- Hardware Independent supports a wide range of diffractometers
- Floating Network License can work as a shareable multi-seat license





### **JADE Pro JADE Pro is all-inclusive!**

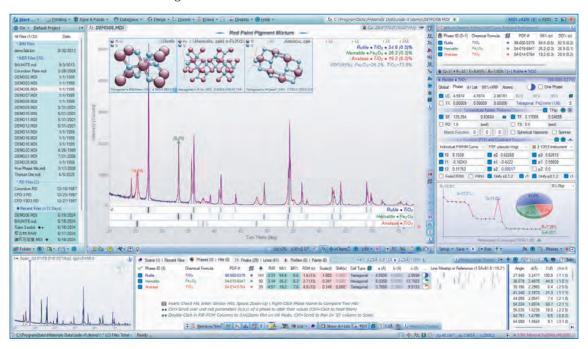
IADE Pro includes all of the features of JADE Standard, plus batch processing Whole Pattern Fitting (WPF) and Rietveld refinement tools that go beyond what is available in JADE Standard.

IADE's One-Click-Analysis  $^{\text{TM}}$  and Toolkit are only found in JADE Pro. Additionally, JADE Pro

features a unique floating network license that allows access on one concurrent system at a time, so when a license is not in use, it is available to someone else in your organization. Discounted additional seats create access to more concurrent shareable licenses that work together.

## **KEY POINTS**

- All-inclusive Everything in JADE Standard and so much more
- Phase ID (Search/Match)
- Batch processing Whole Pattern Fitting (WPF) and Rietveld
- One-Click-Analysis™ for Whole Pattern Fitting (WPF)
- Pattern Indexing (All Crystal Systems)
- Rietveld Structure Refinement (Atomic Parameters)
- Ab Initio Tools (Charge Flipping +)
- Cluster Analysis of Powder Patterns
- Hardware Independent supports a wide range of diffractometers
- Floating Network License can work as a shareable multi-seat license
- Discounted Additional Seats Available





#### **Toolkit is found only in JADE Pro**

**ED-XRF** ▶

For manual or automatic identification of elements and minerals.

#### **XRD Digitizer**

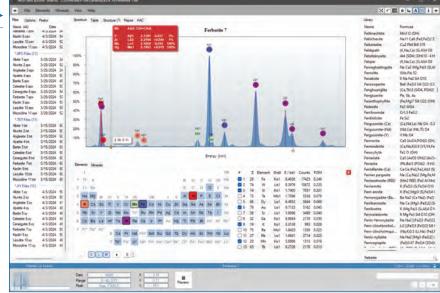
Take a picture of an XRD plot and this tool will convert it to a digital format that you can now work with in JADE. Use advanced image editing options to tweak any imperfections or remove artifacts from your image.

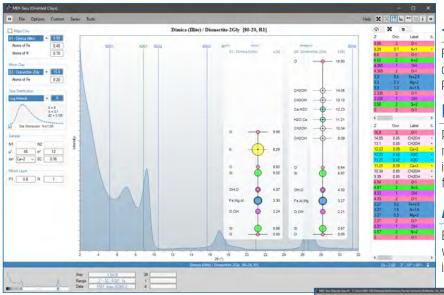
#### Charge Flipper + Solve Structure

For structure solution using the charge flipping method.

#### **Regression Analysis**

Easily perform regression analysis on data without leaving JADE Pro's Toolkit.





#### ◆ For Oriented Clays

For (001)-patterns. Performs simulations and analyses of oriented clays based on structure types introduced by R.C. Reynolds.

#### **Microstructure Tool**

Provides methods for the analysis of the residual stress and microstructure defects for various diffraction geometries and is applicable to cubic phases (FCC, BCC) following a profile fitting in JADE (D. Rafaja).

#### **Area Detector Tool**

Built to extract powder patterns from images obtained with flat area detectors, so that they can easily be analyzed in JADE Pro.

#### Raman Mineral >

S a 1 m3

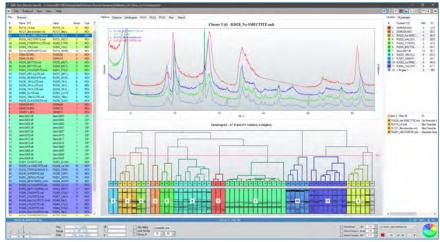
Designed for the mineral phase ID from Raman Stokes spectra. Whole spectra correlation functions are used for finding similarity matches for single-phase and multi-phase spectra.

#### **MDI SizeStrain**

Designed for crystallite size and micro-strain analysis using X-ray powder diffraction data. This Toolkit supports both Warren-Averbach and Williamson-Hall methods for crystallite size determination.

#### **3D Scatter Plot**

A scatter plot tool for xyz data sets from columnar data files.



#### **◄ Cluster Analysis with Dendrograms**

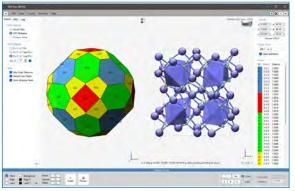
View your clusters with automatic cluster color coding with Principal Component or Fuzzy Clustering. Streamline your analysis by adding reference patterns as cluster markers, recalling setup protocols for reproducibility and other correlation functions for determining pattern similarities.

#### **Pattern Matching**

Try the full Pattern Matching method used in both ROCKJOCK and ClaySim.

#### Laue Visualization

Laue simulations, stereographic projections and alternative structure visualizations.



#### **◄ BFDH Morphology**

Create predicted crystal shapes based on the Bravais-Friedel-Donnay-Harker model. Many options for other calculations.

#### Simulated Annealing ►

An alternative direct space structure solution method (SA)

designed for inorganic structures applying Coulomb interactions and repulsive potentials.



## **JADE Standard**

JADE Standard provides a more limited set of tools than JADE Pro, but without the annual updates and renewal costs. Available in three levels -

- 1. Basic Level
- 2. Add either Phase ID (S/M) or WPF + Rietveld
- 3. Add both Phase ID (S/M) and WPF + Rietveld

Included with all levels of JADE Standard - Simulation and Visualization - Simulate powder patterns to create your own custom database. Visualize a structure in 3D with motion. Visual cues throughout to better help you and your team understand the materials you are analyzing.

## **REY POINTS - ALL LEVELS**

- Reads all vendor pattern files
- Handles variable/fixed-slit data
- Auto peak search and multi-filter search
- Simulate powder patterns from crystal structures
- No annual renewal fee
- Individually licensed for workstations
- Hardware Independent supports a wide range of diffractometers

#### PHASE ID (SEARCH/MATCH)

- Found in Levels 2A: Search/Match 3: Search/Match and Whole Pattern Fitting
- Phase Identification (S/M)
- Minor and Trace Phase ID

#### WPF + RIETVELD

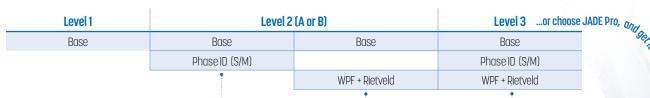
- Found in Levels
  2B: Whole Pattern Fitting (WPF)
  3: Search/Match and WPF
- WPF & Rietveld Refinement











Phase ID (S/M) - JADE Standard Level 2A or Level 3

Profile-based Search/Match with minor and trace phase ID or search on a single peak. Ability to include chemistry filters and unit cell data as well as preferred-orientation.

WPF + RIETVELD - JADE Standard Level 2B or Level 3

Quantify weight percentages and identify minor phases quickly and easily.



## **JADE QA**

A special build meant for production line Quantitative Analysis (QA) typically carried out by a technician. Requires a setup file created by an analytical scientist using JADE Pro. JADE QA runs independently of JADE Pro and was designed to extend the power of Rietveld Quantitative Analysis (QA) to those in your organization responsible for overseeing consistencies in production who may or may not have a background in XRD.





## **Get More Power With JADE Pro**

#### JADE is available in Pro and Standard

1 2A 2B 3

JADE FEATURE LIST	Pro
	PIU
Intuitive, State-of-the-Art UI Reads All Vendor Pattern Files	
	M
Smart Column Data Importer	Y,
Merge Multi-Scans/Segments	Y.
1,2, and 3D Pattern Overlay	
Déjà Vu ID Similar Pattern Files	<b>Y</b> .
Handles Variable/Fixed-Slit Data	Y.
Multiple Display Axes and Any Diffraction Wavelength	ンソンソンソンソンソン
Zero Offset Correction and Theta Calibration	<b>Y</b>
Background Fitting/Batch Fitting/Removal	<b>Y</b>
Auto Peak Search and Multi-Filter Search	<b>Y</b>
Wide Variety of Report and Print Options	<b>V</b>
Includes Starter Database	<b>V</b>
ICDD PDF Database Access (as well as other databases)	<b>~</b>
Multitude of Data Choices for Database Retrieval	<b>V</b>
View Crystal Structures and Reflection Planes	<b>✓</b>
View Polyhedra, Calculate Bond Distances and Angles	<b>✓</b>
Simulate Powder Patterns from Crystal Structures	1
Individual Analyst/Shared Work Environments	1
Support for Custom d-1% list for Phase Overlay	1
Excellent Profile Fitting and Batch Profile Fitting	1
RIR and DDM Quantitative Analysis	ンソンソンソンソン
Size & Strain and Residual Stress Analysis	1
Line-Based Unit Cell and Batch Refinement	1
Structure Database Manager	1
Profile Based Search/Match (S/M)	1
Subfile + Chemistry + Unit Cell Filters in S/M	1
Minor and Trace Phase ID	
Single Peak and Severely-Orientation S/M	
Solid-Solution and Isotypical Phase ID	
Auto-S/M in background thread on loading a pattern file	1
Full pattern S/M on PD3 phases by correlation & similarity	
WPF and Rietveld Refinement (Unlimited Phases)	
Gold Standard Quantitative Analysis	
Seamless Integration with Database Retrieval	
Auto-Loading of Structure Data from Database	
Combination of Phases with/without Structure Data Weight%	
	M
Linking of Unit Cell, Phase Broadening Parameters	ンソンソンソンソン
Custom d-1% (hkl, FWHM, Skew) to Fit Clay Phases	Y,
March-Dollase and Spherical Harmonics Orientation Function	Y,
Amorphous Content with/without Internal Wt% Standards	Y,
Atomic Occupancy and Bond Distance Constraints	<b>Y</b> ,
Polynomial and Spline Curve Background Fitting	<b>V</b>

All levels of JADE Standard are perpetual licenses and sold without the benefit of updates or any of JADE Pro's extensive set of extra

features, like the ability to share a license. Since it is a subscription, JADE Pro has a lower buy-in than JADE Standard.

If you have JADE Classic, there still may be trade-in value for your JADE Classic towards the purchase of a current version of JADE Standard or JADE Pro.

#### Additional JADE Features Only Available in JADE Pro Pro One-Click-Analysis™ - Phase ID and WPF/Rietveld Analysis Size and Strain Analysis (Warren-Averbach & Williamson Hall) Pattern Indexing (All Crystal Systems) Simulate Series of Powder Patterns from Crystal Structures WPF and Rietveld Refinement Batch Processing Rietveld Structure Refinement (Atomic Parameters) Atomic Occupancy and Bond Distance Constraints Fourier Map Tools Ab Initio Tools (Charge Flipping +) Cluster Analysis of Powder Patterns Floating (Shareable) Network License All Updates via Subscription Extract Patterns from Area Detector 2D Images Digitize Any XRD Plot Image Cluster Analysis Visualization Dendrograms Toolkit App Version of ClaySim Toolkit App for Microstructure Analysis Toolkit App for Structure Solver Workflow Laue Simulations and Visualization Toolkit App for Simulated Annealing Toolkit App for Regression Analysis Toolkit App for Pattern Matching Toolkit App for 3D Scatter Plot

#### **Multiple Languages!**

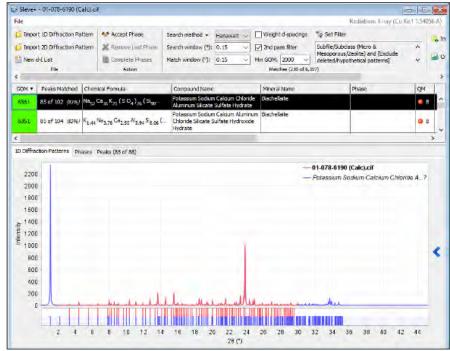
To better support diverse teams sharing a license of JADE Pro across their organization and to ensure consistent results regardless of language, JADE can switch on the fly to Chinese or Japanese languages.



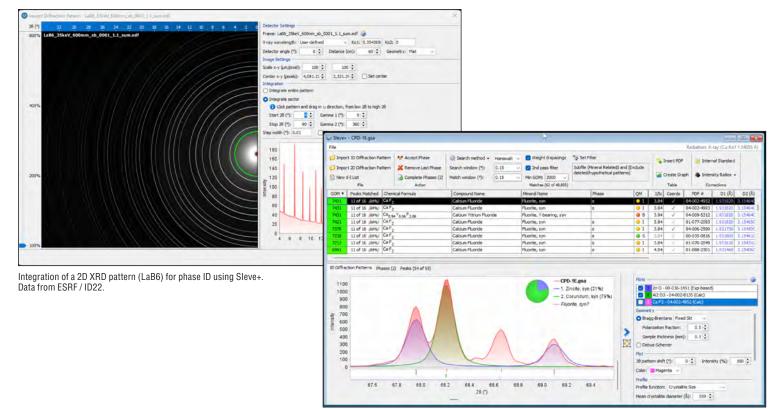
## Sleve/Sleve+ 2026

### Search and Identify

ICDD's search identification programs, Sleve for PDF-2 and Sleve+ for PDF-5+ and PDF-4, are designed to search and identify unknown materials. Sleve/Sleve+ are integrated into the ICDD databases to allow the use of the extensive data mining interfaces, searches, and sorts available to improve the accuracy and precision of the phase identification process.



Phase ID of a zeolite missing the strongest peak using the new Deep Search option in Sleve+.



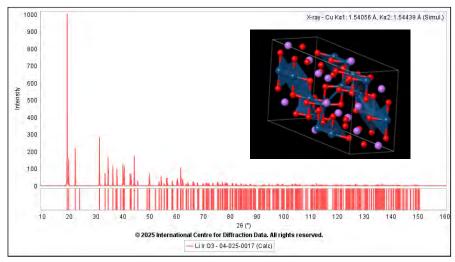
Temporarily hide phase plots with checkboxes in Sleve+.

## **(1)** KEY POINTS

- Included for free in all PDF databases
- Designed to search and identify unknown materials
- Complementary with most commercial software programs
- Match filter algorithms better results, more accuracy
- Identify non-crystalline materials and complex multi-phase specimens
- Strength of Sleve/Sleve+ is evident in the analysis of minor and trace phases
- Supports X-ray and synchrotron data
- "Intelligent" RIR for semi-quantitative analysis\*
  - \* PDF-5+/PDF-4 only feature (Sleve+)

Sleve+ offers a variety of algorithms and options that allow users to optimize results for particular chemistries for both standard and non-standard diffraction systems.





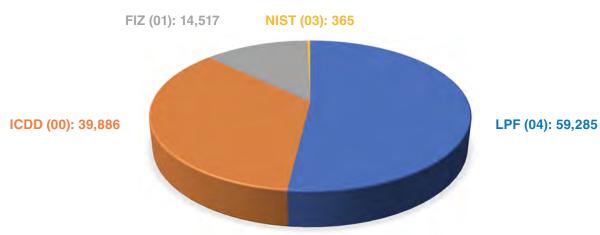
Lithium Iridium Oxide (LiIrO<sub>2</sub>): 04-025-0017

## **KEY POINTS**

- Cost-effective
- Single License: Three (3) year license term
- Optional: Two (2) additional seats (three (3) year license term)
- Six (6) year and nine (9) year prepaid packages available
- New license available after term expires
- Featuring 114,000+ entries, including inorganic and organic entries
- 84,200+ entries with atomic coordinates
- Number of entries with I/I<sub>c</sub> values: 84,800+
- Extensive data mining is facilitated with 118 display fields coupled with 82 searches
- Data focused on ambient entries
- Fundamental mineral entries
- ICDD data mining software Sleve+ Included

#### **Database Sources**

(See inside front cover for ICDD's collaborative partners)



#### **Axiom Purchase Options**

PDF-4/Axiom 2026	USB CODE	DOWNLOAD CODE	2025-2026 LIST \$	2025-2026 Academic \$
New - Single License	AX4S26	AX4L26	\$3,715	\$2,665
Two Additional Seats (each seat: single PC installation)	AX4S26MS	AX4L26MS	\$1,050	\$540
PDF-4/Axiom 2026 Packages*	<u>'</u>	'		
PDF-4/Axiom - 6 years	AX4S6YR	AX4L6YR	\$7,430	\$5,330

AX4S9YR

AX4L9YR

\* Purchase two or three Single Licenses

PDF-4/Axiom - 9 years

- (Release 2026, Release 2029, and Release 2032)
- Each PDF-4/Axiom product is licensed for three (3) years - Shipping fees for all databases billed on the original invoice
- \*Please see product policies on pages 18 and 19

\$11,145

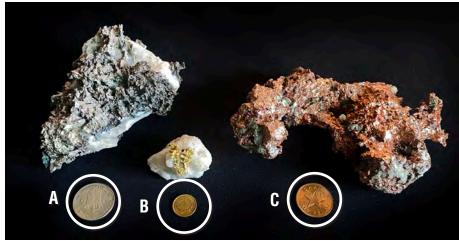
\$7,995



## PDF-4/Minerals 2026

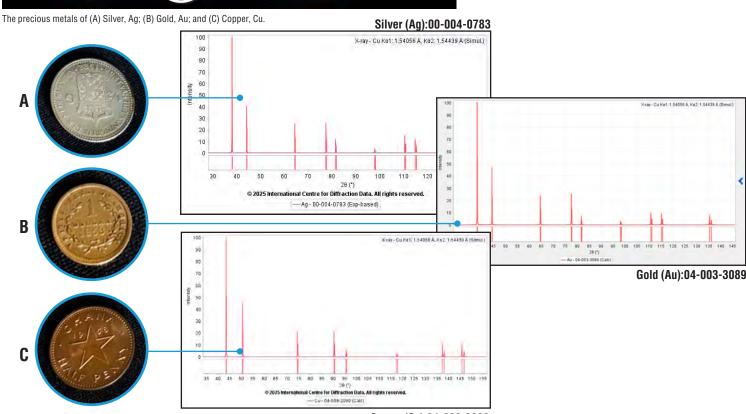
### Comprehensive Mineral Collection

PDF-4/Minerals is the most comprehensive collection of mineral data in the world! Ninety-seven percent of all known mineral types, as defined by the International Mineralogical Association (IMA), are represented in the database.



## **(7)** KEY POINTS

- Rapid phase identification, plus support for automated quantitative analyses
- Features 54,900+ entries, including 45,600+ entries with atomic coordinates
- Integrated data mining software
- All entries have digital patterns for use in total pattern analysis
- More than 43,300+ entries have I/I<sub>c</sub> values for quantitative analysis by Reference Intensity Ratio (RIR)
- Combines powder and single crystal data
- Classified by IMA designations
- A subset of the PDF-5+ database with all the features and capabilities, targeted to minerals and mineral related compounds
- Extensive data mining is facilitated with 118 display fields coupled with 84 searches
- One (1) year license serviced by subscription



#### Copper(Cu):04-009-2090

MNL427MR31

\$4,925

\$3,250

#### **PDF-4/Minerals Purchase Options**

•				
PDF-4/Minerals 2026	USB CODE	DOWNLOAD CODE	2025-2026 LIST \$	2025-2026 Academic \$
New - Single License	MN4S26	MN4L26	\$1,840	\$1,250
Renewal from 2025 to 2026	MNR4S26	MNR4L26	\$985	\$650
PDF-4/Minerals 2026 Multi-year License*				
Renewal - 3 year license	MNS427MR29	MNL427MR29	\$2,955	\$1,950

MNS427MR31

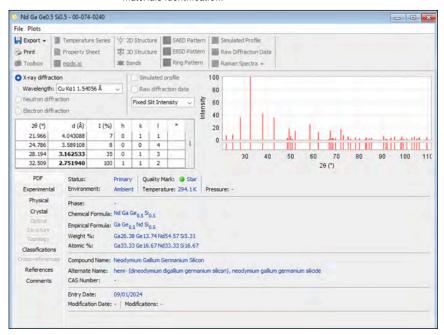
Renewal - 5 year license
\*Please see product policies on pages 18 and 19



## PDF-2 2026

#### Phase Identification + Value

PDF-2 is designed for inorganic materials analyses. In addition, common organic materials from ICDD are added to the database to facilitate rapid materials identification.



## **KEY POINTS**

- Features 369.200+ entries
- 265,700+ entries have I/I<sub>c</sub> values for quantitative analysis by Reference Intensity Ratio (RIR) method
- Integrated data mining software
- Sleve search-indexing software included
- · Combines powder and single-crystal data
- Rapid data mining with our 60 display fields coupled with 75 searches
- · Five (5) year license

Intermetallic compounds display interesting physical and chemical properties. They are used in industry as shape memory alloys, as materials for hydrogen storage, catalysis, as well as magnetic, thermoelectric, or superconducting materials.

#### **PDF-2 Purchase Options**

PDF-2 2026: Five (5) year License	USB CODE	DOWNLOAD CODE	2025-2026 LIST \$	2025-2026 ACADEMIC \$
New - Single License	C2S26PN	C2L26PN	\$9,395	\$6,265
Renewal - from 2025 to 2026	C2S25PU26	C2L25PU26	\$2,115	\$1,375
Renewal - from 2024 to 2026	C2S24PU26	C2L24PU26	\$4,230	\$2,750
Renewal - from 2023 to 2026	C2S23PU26	C2L23PU26	\$6,345	\$4,125
Renewal - from 2022 to 2026	C2S22PU26	C2L22PU26	\$8,460	\$5,500
PDF-2 2026: Ten (10) year License Package				
New - Licenses (2026 and 2031)	PD2S10YR	PD2L10YR	\$18,790	\$12,530
PDF-2 2025 Site License*				
New - Master License	C2S26PNM	C2L26PNM	\$9,395	\$6,265
New - additional licenses, each	C2S26PNC	C2L26PNC	\$2,115	\$1,375
Renewal - Yearly Master License	C2S25PU26M	C2L25PU26M	\$2,115	\$1,375
Renewal - additional licenses, each	C2S25PU26C	C2L25PU26C	\$855	\$590
PDF-2 to PDF-5+ Conversion				
PDF-2 2025 to PDF-5+ 2026	P2525RSU26	P2525RLU26	\$2,115	\$1,375
PDF-2 2024 to PDF-5+ 2026	P2524RSU26	P2524RLU26	\$4,230	\$2,750
PDF-2 2023 to PDF-5+ 2026	P2523RSU26	P2523RLU26	\$6,345	\$4,125
PDF-2 2022 to PDF-5+ 2026	P2522RSU26	P2522RLU26	\$8,460	\$5,500
*Please see product policies on pages 18 and 19	·			

**Every Powder Diffraction File** entry is editorially reviewed and assigned a quality mark.

We are proud to say **ICDD's Quality Management** System is ISO 9001:2015 certified by DEKRA



## Which ICDD Database is Right for You?

	PDF-5+ 2026	PDF-5+ Server Edition 2026	PDF-4/ Axiom 2026	PDF-4/ Minerals 2026	PDF-2 2026
I work primarily with inorganic compounds					
I work primarily with organic compounds					
I work primarily with minerals					
I work with electron diffraction/synchrotron					
I need to be flexible with my work location					
I need to do quantitative analysis					
Rietveld		•	•	•	
Reference Intensity Ratio (RIR)					
Pattern Fitting					
I need a database with atomic coordinates					
How many entries are in the database?	1,126,213	1,126,213	114,053	54,940	369,274
How many entries are new?	25,064	25,064	4,075	1,715	9,867
I am using vendor software for my analysis. Which databases will work with my vendor software?  Global software vendors provide search/match software that interfaces with our databases. Please check with your software vendor for compatibility.	•	•	•	•	•
What is the license term for the database?	1 year	1 year	3 years	1 year	5 years
Multi-year license available					
Site license available		<b>■</b> †			

<sup>†</sup> Requires a 5 concurrent user minimum license.

#### **System Requirements:**

- Intel<sup>®</sup> Core<sup>™</sup> 2 Duo processor (recommended minimum) or better
- 4 GB system memory

#### JADE Standard/JADE Pro:

- Intel® Core™ i7, i5, or i3
- 8 GB system memory
- 24" or larger monitor Your JADE experience will be richer on a larger screen.

#### **Hard Disk Space Requirements**

- PDF-2 2026 5 GB
- PDF-5+ 2026 38 GB
- PDF-5+ Server Edition 2026 38 GB
- PDF-4/Minerals 2026 2 GB
- PDF-4/Axiom 2026 4 GB
- JADE Standard/JADE Pro 400 MB

#### Support Microsoft® Operating System:

- Windows® 7 SP1
- Windows® 8/8.1
- Windows® 10
- Windows® 11

Do you have 5 or more concurrent users on a network? See pages 6 - 7 for Server Edition details.



## **ICDD's FREE Webinar Series** Mastering the PDF Database and JADE Software

Learn the potential of Powder Diffraction File and JADE from ICDD's and Materials Data's experts and developers. Take an hour to build your confidence and productivity.



## **Our Experts Teach You:**

**Quantitative Phase Analysis with Cements** Raman Search Match **Clay Analysis** 

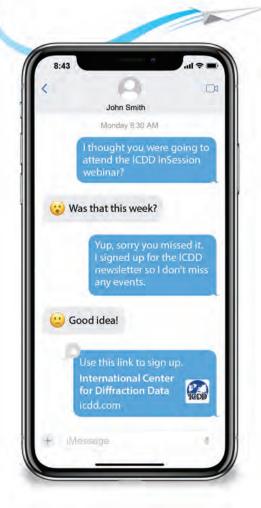
**Data Mining in the Powder Diffraction File Navigating a Powder Diffraction File Entry** and more



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Contact ICDD for PDF-5+ registration information - registration@icdd.com

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#### **RELEASE 2026**

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PDF-4/Axiom	3 years			
PDF-5+*	1 year			
PDF-4/Minerals*	1 year			

<sup>\*</sup> Serviced by subscription featuring automatic annual renewals.

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PDF-5+ The Powder Diffraction Standard



PDF-5+ Server Edition Extend the Power of Your Database -

Built for Teams of Five or More



PDF-4/Axiom

Quality plus Value



PDF-4/Minerals

Comprehensive Mineral Collection



PDF-2

Phase Identification + Value

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\*Dates, location, and format of courses and conferences may be subject to change.

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#### www.icdd.com $\rightarrow$ Conferences

#### 75th Denver X-ray Conference

The Westin Chicago Lombard, Lombard, Illinois, USA 3-7 August 2026

DXC is the world's largest annual conference dedicated to the fields of X-ray analysis. Experts in the fields of X-ray diffraction and X-ray fluorescence gather for a week of workshops, technical sessions, and manufacturer's exhibits. DXC attendees will have access to sessions on the latest advancements in XRD and XRF.

www.dxcicdd.com

## Pharmaceutical Powder X-ray Diffraction Symposium

XRD Training for the Pharmaceutical Scientist! ICDD sponsors this symposium designed to create a forum for the exchange of knowledge and cutting-edge ideas among those interested in the combined fields of XRD and pharmaceutical sciences. Visit the PPXRD website for news on the next symposium. Search and browse presentations from past PPXRD for free.

www.icdd.com/ppxrd

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ICDD's international journal, *Powder Diffraction*, focuses on materials characterization employing X-ray powder diffraction and related techniques. *Powder Diffraction* is publishhed quarterly and with special topical issues. See freshly approved manuscripts online - since 2013, authors have utilized First View to speed approval of their manuscripts.

www.icdd.com/pdj

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mail International Centre for Diffraction Data

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Newtown Square, PA 19073-3273 USA

#### **ICDD California Office**

phone 1.925.449.1084 1.925.373.1659 fax

email mdi@materialsdata.com technical support support@materialsdata.com web www.materialsdata.com

International Centre for Diffraction Data mail

> **DBA Materials Data** 2551 Second Street Livermore, CA 94550 USA

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- Online order tracking
- Email order confirmation
- Email shipping notification
- Customer satisfaction

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Product warranty is on the system requirements associated with the product release. If you need to change your computer specifications, ICDD recommends that you reference the ReadMe.txt or product insert for your product release to check for compatibility.

#### **Payment Terms**

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**NOTE:** Customers in Japan and South Korea, must order through an ICDD licensed distributor. www.icdd.com/distributors

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#### ICDD Pennsylvania:

12 Campus Boulevard
Newtown Square, PA 19073-3273, USA
Tel: 610.325.9814 ◆ Fax: 610.325.9823
Toll Free: 866.378.9331 (US & Canada only)
www.icdd.com



#### ICDD California:

DBA Materials Data 2551 Second Street Livermore, CA 94550, USA Tel: 925.449.1084 ◆ Fax: 925.373.1659 www.materialsdata.com