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Rainer Münz  
Stefan Scherer

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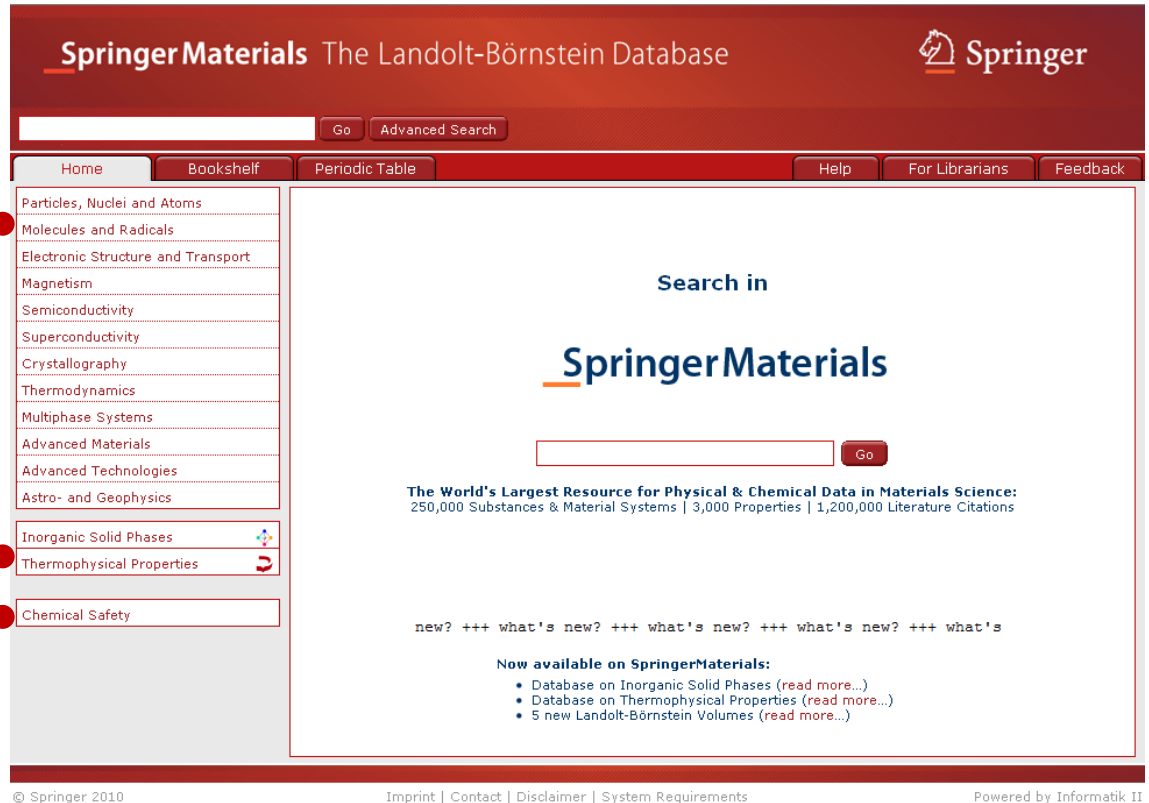
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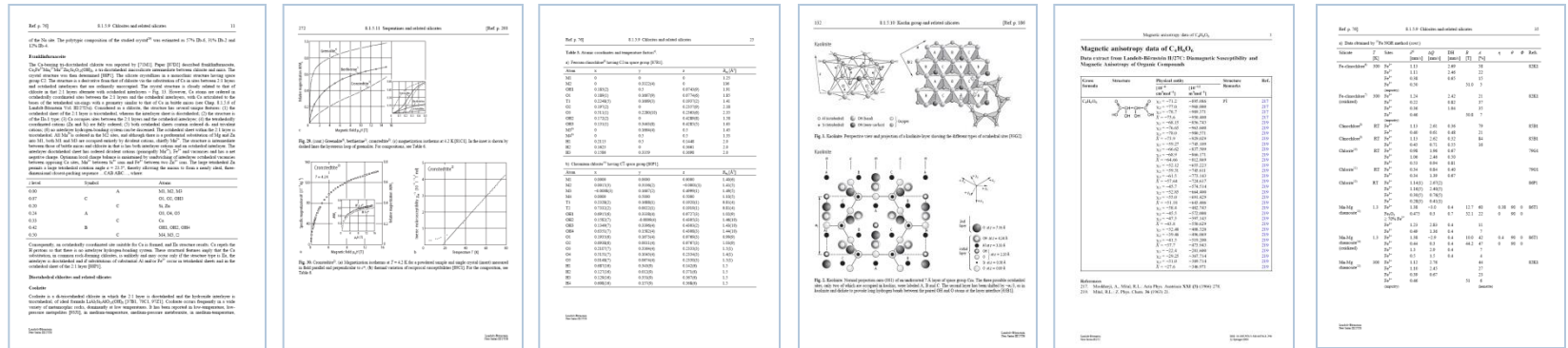
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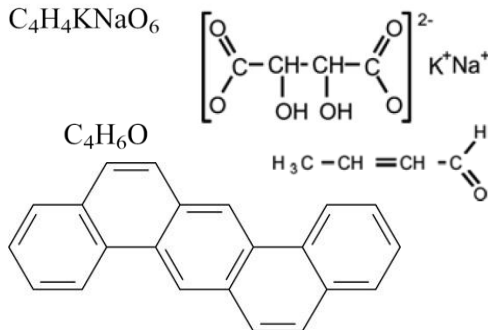
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# What is Landolt-Börnstein?

- > 200,000 pages | > 100,000 online documents | > 150,000 figures



- > 250,000 chemical substances | > 1,200,000 literature references



[94G1]	Goffé, B., Baromet, A., Morin, G.: <i>Eur. J. Mineral.</i> <b>6</b> (1994) 897.	[98B1]	Bailey, S.W.: <i>Clays Clay Miner.</i> <b>36</b> (1988) 193.
[94G2]	Gregori, D.A., Meussler, R.C.: <i>Hyperfine Interact.</i> <b>83</b> (1994) 495.	[98B2]	Bailey, S.W.: <i>Am. Rev. Mineral.</i> <b>39</b> (1988) 347.
[94P1]	Parise, J.B., Leinenweber, K., Weidner, D.J., Tan, K., Von Dreele, R.B.: <i>Am. Mineral.</i> <b>79</b> (1994) 193.	[98B3]	Bernauer, R.G.: <i>J. Petrol.</i> <b>29</b> (1988) 445.
[94R1]	Rancourt, D.G.: <i>Phys. Chem. Miner.</i> <b>21</b> (1994) 250.	[98B4]	Minoura, A., Banno, D., Kawan, F., Fruegy, C.: <i>Phys. Chem. Miner.</i> <b>16</b> (1988) 180.
[94V1]	Villieras, F., Yvon, J., Cases, J.M., De Donato, P., Lhoste, F., Baeza, R.: <i>Clays Clay Miner.</i> <b>42</b> (1994) 679.	[98P1]	Perrot, D.R., Rouse, E.C., Bailey, S.W.: <i>Am. Mineral.</i> <b>73</b> (1988) 876.
[95R1]	Bailey, S.W., Bartsch, J.E., Barker, W.W., Ketchum, G.: <i>Am. Mineral.</i> <b>80</b> (1995) 65.	[98P2]	Rabe, A.C., Radke, F.: <i>Am. Mineral.</i> <b>73</b> (1988) 135.
[95C1]	Catti, M., Ferraris, G., Hall, S., Pavese, A.: <i>Phys. Chem. Miner.</i> <b>22</b> (1995) 200.	[98P3]	Pastarska, H.: <i>Radovity, K., in Proc. 10th Conf. Clay Mineral. Petrol. Ontario, 1988</i> , p. 113.
[95W1]	Welch, M.D., Barras, J., Klinowski, J.: <i>Am. Mineral.</i> <b>80</b> (1995) 441.	[98P4]	Bailey, S.W., Lister, J.S.: <i>Clays Clay Miner.</i> <b>37</b> (1989) 193.
[95Z1]	Zhan, W., Guggenheim, S.: <i>Clays Clay Miner.</i> <b>43</b> (1995) 622.	[99B1]	Benn, A.J., Schreyves, D.: <i>Am. Mineral.</i> <b>74</b> (1989) 1115.
[95Z2]	Zheng, H., Bailey, S.W.: <i>Am. Mineral.</i> <b>80</b> (1995) 387.	[99P1]	Perovskii, J.B., Kodama, H.: <i>Can. Mineral.</i> <b>37</b> (1989) 511.
[96A1]	Andri, M., Wildner, M., Taram, M., Langer, K., Schulz, R.: <i>Phys. Chem. Miner.</i> <b>23</b> (1996) 241.	[99P2]	Zimmerman, V., Frenzel, T., Frenzel, A.M.: <i>Eur. J. Mineral.</i> <b>1</b> (1989) 633.
[96B1]	Baker, J., Holland, T.J.B.: <i>Am. Mineral.</i> <b>81</b> (1996) 676.	[99P3]	Wölsner, D.E.: <i>Am. Mineral.</i> <b>74</b> (1989) 203.
[96G1]	Jullien, M., Baromet, A., Goffé, B.: <i>Am. Mineral.</i> <b>81</b> (1996) 67.		
[97D1]	Droth, M., Trautwein, A.X., König, I., Stuess, E., Bender Koch, C.: <i>Phys. Chem. Miner.</i> <b>24</b> (1997) 281.		
[97G1]	Grevel, K., Fasshauer, D.W., Erzner, S.: <i>Eur. J. Mineral.</i> <b>9</b> (1997) 158.		

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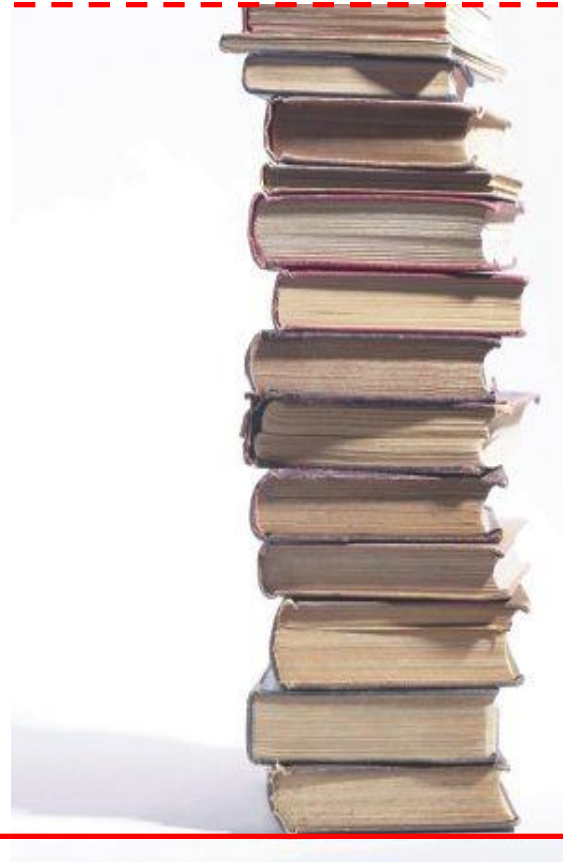


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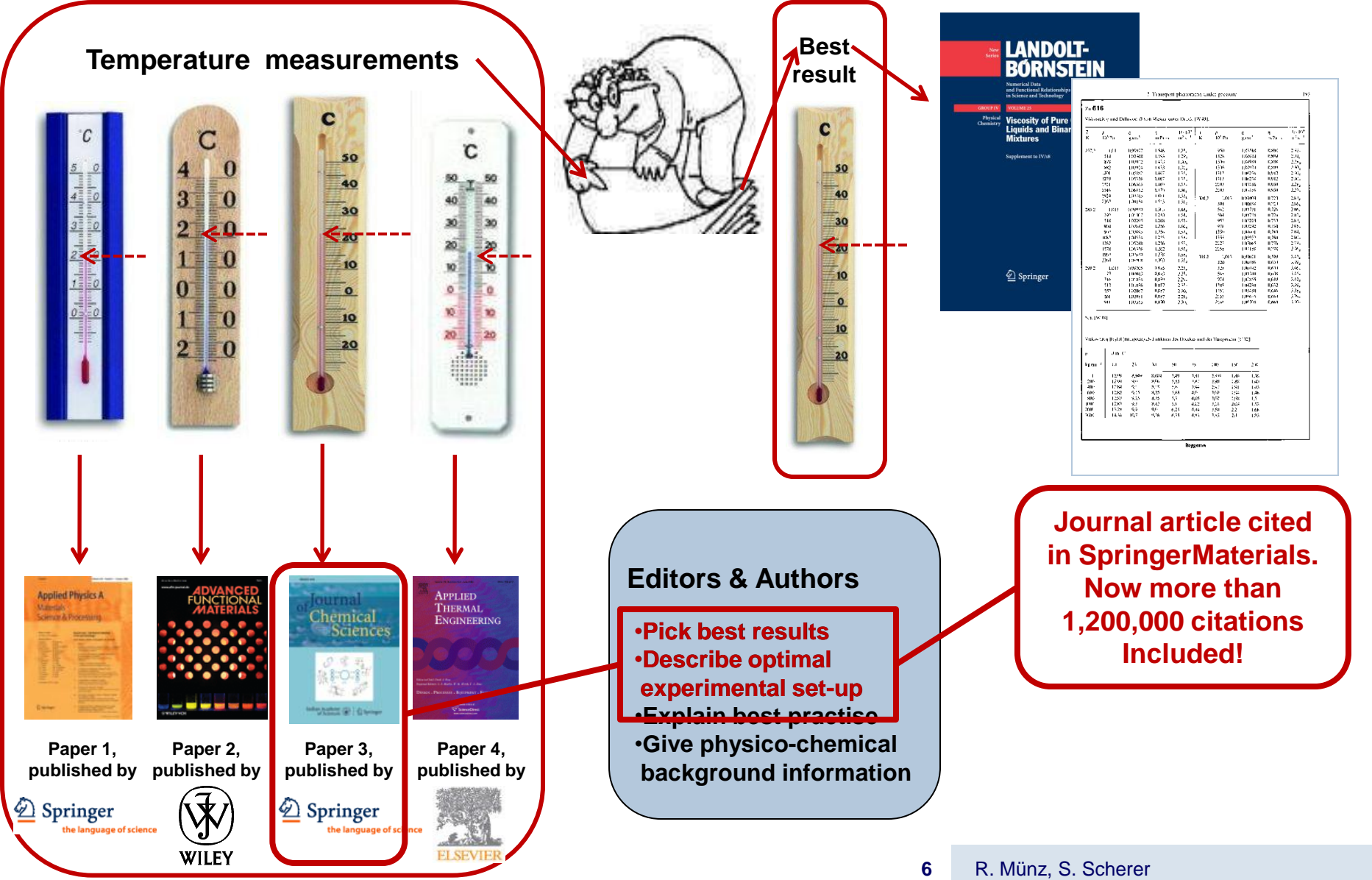


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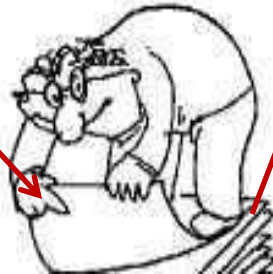
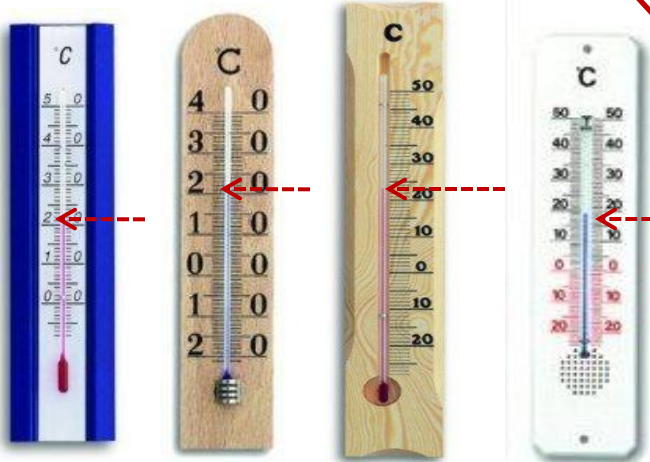
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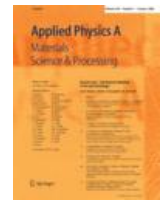
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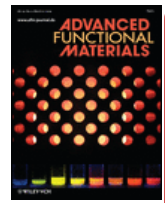
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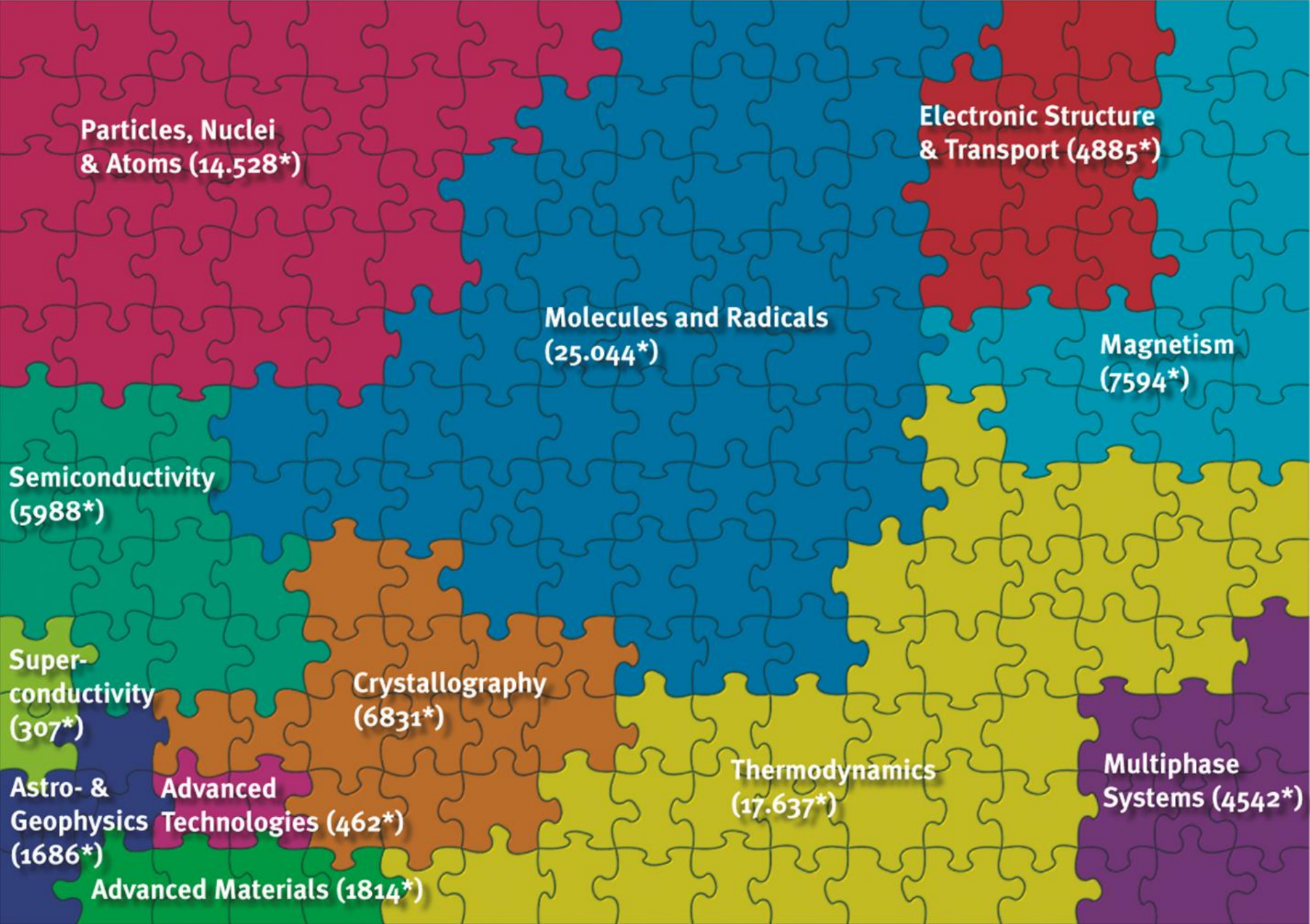
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## Refractory, Hard and Intermetallic Materials

Title Page, Contributors, and Preface  

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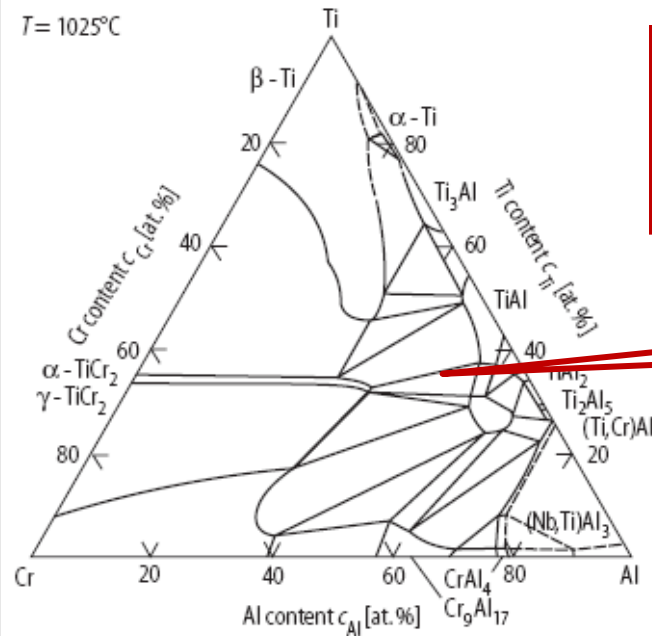
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- Title Page, Authors, Preface
- 1 Introduction (P. Beiss, R. Ruthardt, T. ...)
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- 13 Hard materials (G.E. Spriggs)
- 14 Intermetallic materials (G. Sauthoff)
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    - 14.4.1 Basic properties of constituent intermetallics
    - 14.4.2 Titanium aluminide alloys
      - 14.4.2.1 Phase diagrams
      - 14.4.2.2 Ti(3)Al-base alloys
      - 14.4.2.3 TiAl-base alloys
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    - 14.4.4 Nickel aluminides
    - 14.4.5 Silicides
  - References for 14

T = 1025°C



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Figures

Fig. 14.4. Isothermal section of the Ti-Al-Cr phase diagram at 1025 °C [97P1] - see also [97X1, 98H3].

Text

#### 14.4.2.2 Ti<sub>3</sub>Al-base alloys

##### 14.4.2.2.1 Compositions and applications

Various Ti<sub>3</sub>Al-base alloys have been developed which contain niobium as major alloying element and further elements for obtaining an optimised balance of strength, deformability, toughness and oxidation resistance. The alloys are two-phase or three-phase and contain additional phases besides the phase Ti<sub>3</sub>Al, which is known as the α<sub>2</sub> phase. The usual additional phases are the β phase, which is Ti-rich and which has the b.c.c. structure or is ordered with B2 structure, the orthorhombic Ti<sub>3</sub>Al<sub>2</sub> phase, the Nb-rich phase Ti<sub>3</sub>Al<sub>2</sub>Nb, and/or the ω-type phase Ti<sub>4</sub>Al<sub>3</sub>Nb with B<sub>8</sub> structure [92B1, 95B]. In addition, Ti<sub>3</sub>Al-base alloys have been used as matrices for intermetallic matrix composites, which rely primarily on the presence of the SiC fibre SCS-6 [92L]. Current Ti<sub>3</sub>Al-base alloys with engineering significance are listed in Table 6.

Tables

Table 6. Important Ti<sub>3</sub>Al-base alloys [92B1, 95B].

Type	Alloy [at.%]	Designation	Phases
α <sub>2</sub>	Ti-24Al-11Nb	24-11	α <sub>2</sub> + β
α <sub>2</sub>	Ti-25Al-11Nb	25-11	α <sub>2</sub> + β

**Bibliographic Information for 10.1007/10858641\_18**

**Cite as:**

Sauthoff, G.: *14.4 Intermetallic materials for structural high temperature applications*. Beiss, P., Ruthardt, R., Warlimont, H. (ed.). SpringerMaterials - The Landolt-Börnstein Database (<http://www.springermaterials.com>). DOI: 10.1007/10858641\_18

**Source:**

<b>Title</b>	<b>14.4 Intermetallic materials for structural high temperature applications</b>
<b>In</b>	14 Intermetallic materials
<b>Author</b>	G. Sauthoff
<b>Part of</b>	Landolt-Börnstein - Group VIII Advanced Materials and Technologies Numerical Data and Functional Relationships in Science and Technology
<b>Volume</b>	<b>2A2:</b> Powder Metallurgy Data. Refractory, Hard and Intermetallic Materials
<b>Edited by</b>	P. Beiss, R. Ruthardt, H. Warlimont
<b>Chapter-DOI</b>	10.1007/10858641_18
<b>Book-DOI</b>	10.1007/b83029 (Volume in Bookshelf)

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- Figs. 13.1 - 13.91

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  - Basic properties of constituent phases
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    - Phase diagrams
    - Ti<sub>3</sub>Al-base alloys
    - TiAl-base alloys
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    - Fe<sub>3</sub>Al-base alloys

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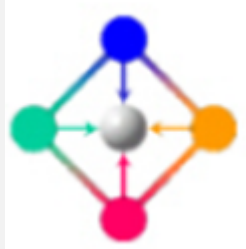
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## Database on Inorganic Solid

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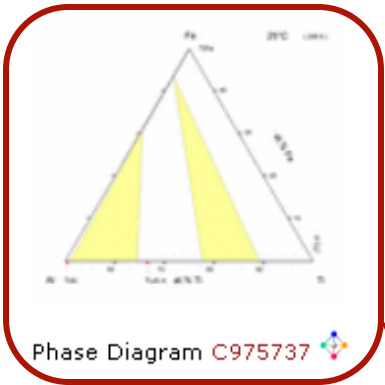
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- **Phase Diagrams**  
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**Al-Fe-Ti, ternary phase diagram, isothermal section**

Element System: Al-Fe-Ti

**Inorganic Solid Phases · phase diagrams**

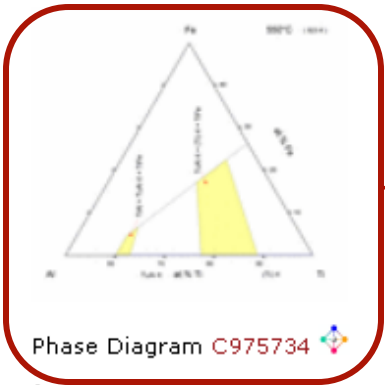
Diagram details: Al conc.[0-50 at.%] vs. Fe conc.[0-50 at.%] vs. Ti conc.[50-100 at.%]



Phase Diagram C975737

from:  
Seibold A., Seibold A., Seibold A., Seibold A.: "Phasengleichgewichte in den ternären Systemen Ti-Fe-O und Ti-Al-Fe", Z. Metallkd. **72** (1981) 712-719.

Diagram details: Al conc.[0-30 wt.%] vs. Fe conc.[0-30 wt.%] vs. Ti conc.[70-100 wt.%]



Phase Diagram C975734

from:  
Volkova M.A., Kornilov I.I.: "PHASE EQUILIBRIUM IN TITANIUM-RICH Ti-Al-Fe ALLOYS", Russ. Metall. **3** (1970) 134-137.

**„Inorganic Solid Phases“:  
Overview of Relevant  
Data Sheets**

Diagram details: Al conc.[0-45 at.%] vs. Fe conc.[0-27 at.%] vs. Ti conc.[55-100 at.%]

Element System: Al-Fe-Ti

Ternary Phase Diagram

Investigated Concentration Ranges: Al conc.[0-50 at.%) vs. Fe conc.[0-50 at.%) vs. Ti conc.[50-100 at.%)

**Constitution Entry Summary**

<b>System</b>	Al-Fe-Ti
<b>Diagram Type</b>	ternary, isothermal section
<b>Concentration Range</b>	partial composition; Ti-TiAl-TiFe / 50-100 at.% Ti
<b>Temperature</b>	298 K
<b>Investigation</b>	experimental
<b>APDIC</b>	non-APDIC

**Reference**

Seibold A.: "Phasengleichgewichte in den ternären Systemen Ti-Fe-O und Ti-Al-Fe", Z. Metallkd. 72 (1981) 712-719.

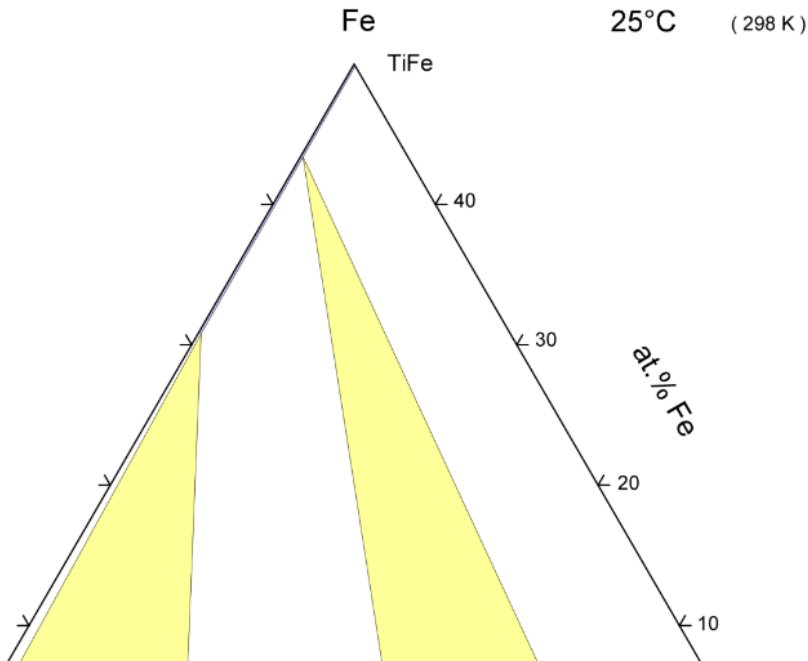
**Phase Diagram**

**Published Phase Diagram**

<b>Figure</b>	19
<b>Title</b>	Al-Fe-Ti isothermal section at 25 °C
<b>Parameters</b>	Al conc.[0-50 at.%) vs. Fe conc.[0-50 at.%) vs. Ti conc.[50-100 at.%)

**Redrawn Phase Diagram**

<b>Concentration range</b>	partial composition; Ti-TiAl-TiFe
<b>Enlargement</b>	0.6



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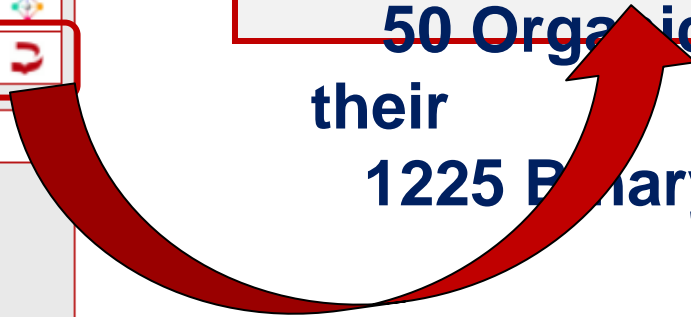
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Subset of the DDBST /  
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Software

& Separation Technology

- Thermophysical Properties of  
50 Organic Liquids + Water and  
their  
1225 Binary Mixtures



**Methanol / Water**

**Thermophysical Data in the Dortmund Data Bank**


**Components**

No.	Formula	Molar Mass	CAS Registry Number	Name
1	CH <sub>4</sub> O	32.042	67-56-1	Methanol
2	H <sub>2</sub> O	18.015	7732-18-5	Water

**List of Available Properties**

- Activity Coefficients at Infinite Dilution ↻
- Azeotropic Data ↻
- Excess Heat Capacities ↻
- Heat Capacities ↻
- Excess Enthalpies ↻
- Solid-Liquid Equilibria ↻
- Excess Volumes ↻
- Densities ↻
- Vapor-Liquid Equilibria ↻

**Thermophysical Properties“: Overview of Available Data**

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**Methanol / Water**  
**Activity Coefficients at Infinite Dilution**
**Components**

No.	Formula	Molar Mass	CAS Registry Number	Name
1	CH <sub>4</sub> O	32.042	67-56-1	Methanol
2	H <sub>2</sub> O	18.015	7732-18-5	Water

**Water solved in Methanol**

Temperature [K]	Activity Coeff. at Inf. Dil.	Measurement Method	Source
298.15	1.6300	CPEQ	4
298.15	4.1800	OTHR	2
307.75	1.6000	EBUL	1
313.15	2.2000	STAT	3
317.85	1.8000	EBUL	1
323.15	3.2600	OTHR	2
327.85	1.7000	EBUL	1
333.15	1.7700	CPEQ	4
337.05	1.7600	EBUL	1
373.15	1.8700	CPEQ	4

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**List of measurement methods**

Short Term	Explanation
CPEQ	calculated from phase equilibria
EBUL	ebullimetry
OTHR	other techniques
STAT	static method

**List of References**

1	Bergmann D.L., Eckert C.A.: Measurement of Limiting Activity Coefficients for Aqueous Systems by Differential Ebullimetry. Fluid Phase Equilib. 63 (1991) 141-150
2	Schmidt T.W.: Determination of Infinite Dilution Activity Coefficients (gamma infinite) using Molecular Beams. US-Patent (1980) 1-16
3	Belousov V.P., Makarova N.I., Sabylin I.I.: Thermodynamic Properties of the Water-Ethanol-Dioxan Ternary System at 25°C. Viniti (1973) 1-10
4	Pierotti G.J., Deal C.H., Derr E.L.: (Addendum included). Ind.Eng.Chem. 51 (1959) 95-102

**Methanol solved in Water**

Temperature [K]	Activity Coeff. at Inf. Dil.	Measurement Method	Source
273.35	1.2300	STAT	26
283.15	1.4000	STAT	26
283.15	1.4100	DILU	22
293.15	1.5500	DILU	22
293.15	1.5500	STAT	26
293.15	1.6800	STAT	16
293.15	2.6900	GLCI	1
297.45	2.1200	GLCI	2
298.15	1.5300	HENR	15
298.15	1.5300	CPEQ	23
298.15	1.5800	GLCI	17
298.15	1.6400	GLCR	8
298.15	1.6400	DILU	6

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- **REACH** (Registration, Evaluation, Authorization and Restriction of Chemicals)
- **GHS** (Globally Harmonized System)
- **RoHS** (Restriction of Hazardous Substances)
- **WEEE** (Waste from Electrical and Electronic Equipment)


- 44,000 documents on **REACH | GHS | RoHS | WEEE**
- Identifies recognized potential dangers of a substance
- Standardized classification and labeling
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European regulations regarding benzenemethanol (C<sub>7</sub>H<sub>8</sub>O)


**Name** benzenemethanol      **Formula:** C<sub>7</sub>H<sub>8</sub>O  
**CAS-RN** 100-51-6      **Molecular Weight:** 108.138 g/mol  
**EG-Index:** 603-057-00-5 (93/72/EEC)  
**EINECS:** 202-859-9 (EINECS)

## Hazard Information (Dangerous Substances Directive 67/548/EEC)

<b>Hazard symbol</b>		93/72/EEC
<b>R-Phrase</b>	Xn Harmful 20/22 R20/22 Harmful by inhalation and if swallowed.	93/72/EEC
<b>S-Phrase</b>	(2-)26 S2 Keep out of the reach of children. S26 In case of contact with eyes, rinse immediately with plenty of water and seek medical advice.	93/72/EEC

## GHS classification (Globally Harmonized System)

Regulation on Classification, Labelling and Packaging of Substances and Mixtures (CLP)

<b>Signal Word</b>	Warning	EC/1272/2008
<b>Pictogram</b>		
<b>Hazard Statements</b>	H332 Harmful if inhaled. H302 Harmful if swallowed.	
<b>Precautionary Statements</b>	P261 Avoid breathing dust/fume/gas/mist/vapours/spray. P264 Wash ... thoroughly after handling. P270 Do not eat, drink or smoke when using this product. P271 Use only outdoors or in a well-ventilated area. P301+P312 IF SWALLOWED: Call a POISON CENTER or doctor/physician if you feel unwell. P304+P340 IF INHALED: Remove victim to fresh air and keep at rest in a position comfortable for breathing. P312 Call a POISON CENTER or doctor/physician if you feel unwell. P330 Rinse mouth. P501 Dispose of contents/container to ...	
<b>Classification</b>	Acute Tox. 4 *	
<b>EC-Reference</b>	Acute Tox. 4 * 603-057-00-5	

Actual Safety Document

## European Chemicals Agency (ECHA) pre-registration

<b>Pre-registered substance</b>	EC-Number: 202-859-9 CAS-Number: 100-51-6 Name: benzyl alcohol Synonym: - Registration Date: 30/11/2010	ECHA pre
---------------------------------	---	----------

## References

**93/72/EEC**  
**Short:** 93/72/EEC  
**Title:** Commission Directive 93/72/EEC of 1 September 1993 adapting to technical progress for the nineteenth time Council Directive 67/548/EEC on the approximation of the laws, regulations and administrative provisions relating to the classification, packaging and labelling of dangerous substances  
**Author:** Council of the European Economic Community  
**Source:** *Official Journal of the European Communities*  
**Volume:** L 258 A (16.10.1993)  
**Page:** 1-1409  
**Year:** 1993  
**Keyword:** hazardous materials; hazard classification  
**Internet Resource:** <http://eur-lex.europa.eu/LexUriServ/LexUriServ.do?uri=CELEX:31993L0072:EN:HTML>  
**Publish Date:** 1993/10/16  
**Short:** EC/1272/2008  
**Title:** REGULATION (EC) No 1272/2008 OF THE EUROPEAN PARLIAMENT AND OF THE COUNCIL of 16 December 2008 on classification, labelling and packaging of substances and mixtures, amending and repealing Directives 67/548/EEC and 1999/45/EC, and amending Regulation (EC) No 1907/2006  
**Author:** European Parliament; Council of the European Union  
**Source:** *Official Journal of the European Union*

EC/1272/2008

# Summary

**LANDOLT-  
BÖRNSTEIN**

**“Landolt-Börnstein”**

ca. 100,000 documents



**Database on Inorganic Solid Phases**

“Linus Pauling Files”

ca. 190,000 documents



**Database on Thermophysical Properties**

“Dortmund Data Bank Software & Separation Technology”

ca. 300,000 data points

**Chemical Safety Data**

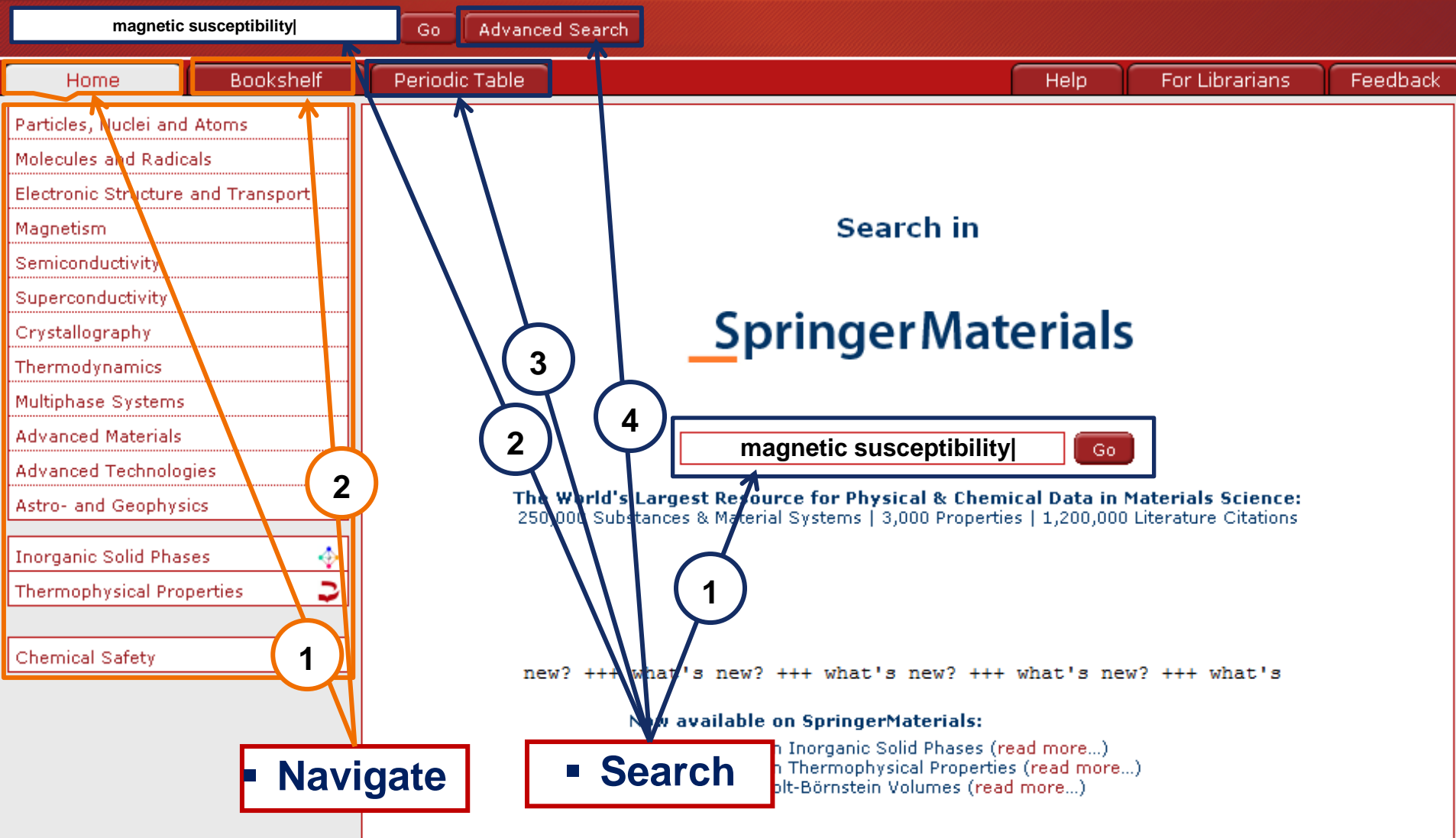
REACH | GHS | RoHS |  
WEEE



ca. 44,000 documents

- Researcher is given the task to replace thick zinc coating with a zinc-manganese coating to prevent corrosion on car bodies
  - Formulated theory on best composition of the compound
  - After testing, result opposite of what was expected
    - Searched for phase diagram to find out why
    - **After 6 weeks of searching** found a Xerox copy of an original article published in an obscure German journal from **1948!**
  - More questions followed as to the crystallography of the compound
  - 2 more weeks of searching in another source to find answer
- With SpringerMaterials :
  - **30 seconds to search and find!**

# Springer Materials



The screenshot shows the SpringerMaterials website interface. At the top, there is a search bar containing the text "magnetic susceptibility" and a "Go" button. Below the search bar, there are navigation tabs: "Home", "Bookshelf", "Periodic Table", "Help", "For Librarians", and "Feedback". On the left side, there is a vertical menu with various categories such as "Particles, Nuclei and Atoms", "Molecules and Radicals", "Electronic Structure and Transport", "Magnetism", "Semiconductivity", "Superconductivity", "Crystallography", "Thermodynamics", "Multiphase Systems", "Advanced Materials", "Advanced Technologies", "Astro- and Geophysics", "Inorganic Solid Phases", "Thermophysical Properties", and "Chemical Safety".

Annotations on the page include:

- A red box labeled "Navigate" with an arrow pointing to the "Bookshelf" tab.
- A red box labeled "Search" with an arrow pointing to the search bar.
- Four numbered circles (1, 2, 3, 4) with arrows pointing to specific elements on the page: circle 1 points to "Chemical Safety", circle 2 points to "Bookshelf", circle 3 points to "Periodic Table", and circle 4 points to the search bar.

The main content area displays "Search in SpringerMaterials" and "The World's Largest Resource for Physical & Chemical Data in Materials Science: 250,000 Substances & Material Systems | 3,000 Properties | 1,200,000 Literature Citations". Below this, there is a search bar with "magnetic susceptibility" and a "Go" button. Further down, there is a section titled "New available on SpringerMaterials:" with links to "Inorganic Solid Phases (read more...)", "Thermophysical Properties (read more...)", and "Landolt-Börnstein Volumes (read more...)".



# Navigate:

- **Subject Areas**
- **Bookshelf**

Go

Advanced Search

Home

Bookshelf

Periodic Table

Help

For Librarians

Feedback

- Particles, Nuclei and Atoms
- Molecules and Radicals
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- Crystallography
- Thermodynamics
- Multiphase Systems
- Advanced Materials
- Advanced Technologies**
- Astro- and Geophysics

Inorganic Solid Phases



Thermophysical Properties



Chemical Safety

Search in

# SpringerMaterials

Go

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250,000 Substances & Material Systems | 3,000 Properties | 1,200,000 Literature Citations

new? +++ what's new? +++ what's new? +++ what's new? +++ what's

Now available on SpringerMaterials:

- Database on Inorganic Solid Phases (read more...)
- Database on Thermophysical Properties (read more...)
- 5 new Landolt-Börnstein Volumes (read more...)

▪ **Select Subject Area**





Go Advanced Search

Home Bookshelf Periodic Table Help For Librarians Feedback

Particles, Nuclei and Atoms  
Molecules and Radicals  
Electronic Structure and Transport  
Magnetism  
Semiconductivity  
Superconductivity  
Crystallography  
Thermodynamics  
Multiphase Systems  
Advanced Materials  
**Advanced Technologies**  
Astro- and Geophysics

> Advanced Technologies

### Advanced Technologies

- Energy
  - Fossil Energy
  - Nuclear Energy
  - Renewable Energy
- Laser
  - Laser Fundamentals
  - Laser Systems
  - Laser Applications
- Metallurgy
  - Powder Metallurgy
  - Metals and Magnets**
  - Refractory, Hard and Intermetallic Materials**
  - Creeping and Rupture Behavior
- Metal Forming Data
  - Ferrous Alloys
    - Index of steel grades  
    - Introduction and Definitions
    - Origin, Preparation and Selection of Data and References  
    - Data on Deformation Behaviour

▪ Select Content



▪ Select Subject Area



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Home Bookshelf Periodic Table Help For Librarians Feedback

> Advanced Technologies



## Refractory, Hard and Intermetallic Materials



Title Page, Contributors, and Preface  



Introduction  



Refractory metals  



Hard materials



Introduction  



Raw materials used in the production of hard materials  



Production methods  

Properties of hardmetals and cermets  



Properties of diamond and cubic boron nitride  



Applications  



Health, safety and ecology  



Acknowledgements  



Figures



Figures 1  

Figures 2  



Figures 3  



Figures 4  



Figures 5  



References  



Intermetallic materials

Introduction  

Topical intermetallic materials  

Production methods  

Intermetallic materials for structural high temperature applications  

References  

**Bread-crum Trail**

**Select Document**

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Help

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Feedback

- Particles, Nuclei and Atoms
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- Astro- and Geophysics

- Inorganic Solid Phases
- Thermophysical Properties

Chemical Safety

> Advanced Technologies

### Refractory, Hard and Intermetallic Materials

- Title Page, Contributors, and Preface
- Introduction
- Refractory metals
- Hard materials
  - Introduction
  - Raw materials used in the production of hard materials
  - Production methods
  - Properties of hardmetals and cermets
  - Properties of diamond and cubic boron nitride
  - Applications
  - Health, safety and ecology
  - Acknowledgements
  - Figures
    - Figures 1
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    - Figures 4
    - Figures 5
  - References
- Intermetallic materials
  - Introduction
  - Topical intermetallic materials
  - Production methods
  - Intermetallic materials for structural high temperature applications
  - References

▪ Select InfoDoc



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Home Bookshelf Periodic Table Help For Librarians Feedback

Group I	Elementary Particles, Nuclei and Atoms
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Group VI	Astronomy and Astrophysics
Group VII	Biophysics
Group VIII	Advanced Materials and Technologies

Group VIII: Advanced Materials and Technologies

**Group VIII: Advanced Materials and Technologies**

- VIII/1a1 Advanced Materials and Technologies · Laser Physics and Applications · Laser Fundamentals, Part 1
- VIII/1a2 Advanced Materials and Technologies · Laser Physics and Applications · Laser Fundamentals, Part 2
- VIII/1b1 Advanced Materials and Technologies · Laser Physics and Applications · Laser Systems, Part 1
- VIII/1b2 **NEW** Advanced Materials and Technologies · Laser Physics and Applications · Laser Systems, Part 2
- VIII/1c Advanced Materials and Technologies · Laser Physics and Applications · Laser Applications
- VIII/2a1 Advanced Materials · Materials · Powder Metallurgy Data, Metals and Magnets
- VIII/2a2 Advanced Materials · Materials · Powder Metallurgy Data, Refractory, Hard and Intermetallic Materials
- VIII/2b Advanced Materials · Materials · Creep Properties of Heat Resistant Steels and Superalloys
- VIII/2c1 **NEW** Advanced Materials · Materials · Metal Forming Data, Ferrous Alloys
- VIII/3a Advanced Materials and Technologies · Energy Technologies · Fossil Energy
- VIII/3b Advanced Materials and Technologies · Energy Technologies · Nuclear Energy
- VIII/3c Advanced Materials and Technologies · Energy Technologies · Renewable Energy
- VIII/4 Advanced Materials and Technologies · Radiological Protection
- VIII/5a Advanced Materials · Physical Properties of Liquid Crystals · Subvolume A
- VIII/2c1 **NEW** Advanced Materials and Technologies · Polymers · Polymer Solutions, Physical Properties and their Relations I (Thermodynamic Properties: Phase Equilibria)

**Navigation via Bookshelf**

**Volumes, Sub-Volumes as in Print**

**Pick Volume**

Go Advanced Search

Home Bookshelf Periodic Table Help For Librarians Feedback

Group I	Elementary Particles, Nuclei and Atoms
Group II	Molecules and Radicals
Group III	Condensed Matter
Group IV	Physical Chemistry
Group V	Geophysics
Group VI	Astronomy and Astrophysics
Group VII	Biophysics
Group VIII	Advanced Materials and Technologies

> Group VIII: Advanced Materials and Technologies

**Advanced Materials · Materials · Powder Metallurgy Data. Refractory, Hard and Intermetallic Materials**  
**Volume: 2A2 · DOI: 10.1007/b83029 · ISBN: 978-3-540-42961-6 · Publisher: Springer-Verlag · Copyright: 2002**

- Title Page, Authors, Preface
- Title Page
- Editors, Authors
- Preface
- SEM Photograph
- 1: Introduction (P. Beiss, R. Ruthardt, H. Warlimont)
- 12: Refractory metals (G. Leichtfried)
  - 12.1: Introduction
  - 12.2: Powder production
  - 12.3: Powder compaction
  - 12.4: Sintering
  - 12.5: Properties
  - References for 12
- 13: Hard materials (G.E. Spriggs)
- 14: Intermetallic materials (G. Sauthoff)
  - 14.1: Introduction
  - 14.2: Topical intermetallic materials
  - 14.3: Production methods
  - 14.4: Intermetallic materials for structural high temperature applications
    - 14.4.1: Basic properties of constituent phases
    - 14.4.2: Titanium aluminide alloys
    - 14.4.3: Iron aluminides
    - 14.4.4: Nickel aluminides
    - 14.4.5: Silicides
  - References for 14

**Browse Table of Contents**

14.4.4: Nickel aluminides

# Search:

- **Google-like**
- **Periodic Table**
- **Advanced Search**



Al-Fe-Ti

Home Bookshelf Periodic Table Help For Librarians Feedback

- 0 Particles, Nuclei and Atoms
- 0 Molecules and Radicals
- 0 Electronic Structure and Transport
- 6 Magnetism
- 0 Semiconductivity
- 1 Superconductivity
- 1 Crystallography
- 0 Thermodynamics
- 5 Multiphase Systems
- 0 Advanced Materials
- 0 Advanced Technologies
- 0 Astro- and Geophysics

9 Inorganic Solid Phases

0 Thermophysical Properties

**Number of Documents Found**

Results 1 - 10 of 22 Documents

previous 1 2 3 next

Multiphase Systems > Ternary Alloys > Phase Diagrams, Crystallography and Thermodynamics > Light Metal Systems > Aluminum (Al-X-Y) Ternary Alloys

Al-Fe-Ti

**Metadata - Substance: Al-Fe-Ti ... Metadata - Element System: Al-Fe-Ti ... Fulltext:** Equilibrium Diagram of the Al-Fe-Ti System and the Segregation of ... Equilibrium Diagram of the Al-Fe-Ti System" (in Russian), Izv. ... #, \*, 12) Raghavan, V., "The Al-Fe-Ti (Aluminum-Iron-Titanium) ... Prop., 40) Raghavan, V., "Al-Fe-Ti (Aluminum-Iron-Titanium).", ...

Multiphase Systems > Ternary Alloys > Phase Diagrams, Crystallography and Thermodynamics > Iron Systems > Selected Systems from Al-B-Fe to C-Co-Fe

Aluminium - Iron - Titanium

**Metadata - Substance: Al-Fe-Ti ... Metadata - Element System: Al-Fe-Ti ... Fulltext:** 1. Investigations of the Al-Fe-Ti Phase Relations, Structures ... points [ 1982All ]. Fig. 1. Al-Fe-Ti. Vertical section at a ... Al content of 25 at.% Fig. 2. Al-Fe-Ti. Vertical section at a ... summarized in Fig. 3. Fig. 3. Al-Fe-Ti. Vertical section at a ...

Inorganic Solid Phases

Al-Fe-Ti, ternary phase diagram, isothermal section

**Metadata - Element System: Al-Fe-Ti ... Fulltext:** Al-Fe-Ti, ternary phase diagram, ... section Element System: Al-Fe-Ti Inorganic Solid Phases · ... from: Raghavan V.: "The Al-Fe-Ti (Aluminum-Iron-Titanium) ... from: Raghavan V.: "The Al-Fe-Ti (Aluminum-Iron-Titanium) ... from: Raghavan V.: "The Al-Fe-Ti (Aluminum-Iron-Titanium) ...

Inorganic Solid Phases

Al-Fe-Ti, ternary phase diagram, liquidus projection

**Metadata - Element System: Al-Fe-Ti ... Fulltext:** Al-Fe-Ti, ternary phase diagram, ... projection Element System: Al-Fe-Ti Inorganic Solid Phases · ... from: Raghavan V.: "The Al-Fe-Ti (Aluminum-Iron-Titanium) ...

Inorganic Solid Phases

Al-Fe-Ti, ternary phase diagram, isothermal section

**Metadata - Element System: Al-Fe-Ti ... Fulltext:** Al-Fe-Ti, ternary phase diagram, ... section Element System: Al-Fe-Ti Inorganic Solid Phases · ...

Inorganic Solid Phases

TiFeAl, crystallographic data

**Metadata - Element System: Al-Fe-Ti ... Fulltext:** data Element System: Al-Fe-Ti Inorganic Solid Phases · ...

**Source of Documents Found**

Suchen nach:

**al-fe-ti in das aktuelle PDF-Dokument**

Ergebnisse:

**1 Dokument(e) mit 67 Treffer(n)**

Neue Suche

Ergebnisse:

- Al-Fe-Ti. Partial isothermal section at 1
- Al-Fe-Ti 27 Fig. 11. Al-Fe-Ti. Partial is
- Al-Fe-Ti. Partial isothermal section at 1
- Al-Fe-Ti Fig. 12. Al-Fe-Ti. Isothermal
- Al-Fe-Ti. Isothermal section at 1000°C
- Al-Fe-Ti. Isothermal section at 900°C.
- Al-Fe-Ti 29 Landolt-Börnstein New Se**
- Al-Fe-Ti Fig. 14. Al-Fe-Ti. Isothermal
- Al-Fe-Ti. Isothermal section at 800°C
- Al-Fe-Ti 31 Fig. 15. Al-Fe-Ti. Partial is
- Al-Fe-Ti. Partial isothermal section at 5
- Al-Fe-Ti References [1940Nis] Nishim
- the Al-Fe-Ti System and the Segregat
- the Al-Fe-Ti System" (in Russian), Izv.
- Al-Fe-Ti 33 Landolt-Börnstein New Se
- Al-Fe-Ti [1987Rag] Raghavan, V., "T
- The Al-Fe-Ti (Aluminum-Iron-Titanium)
- Al-Fe-Ti 35 [1992Gho] Ghosh, G., "Al
- Al-Fe-Ti (Aluminum-Iron-Titanium)", J.
- Al-Fe-Ti [1996Mar] Marchon, I., Saut

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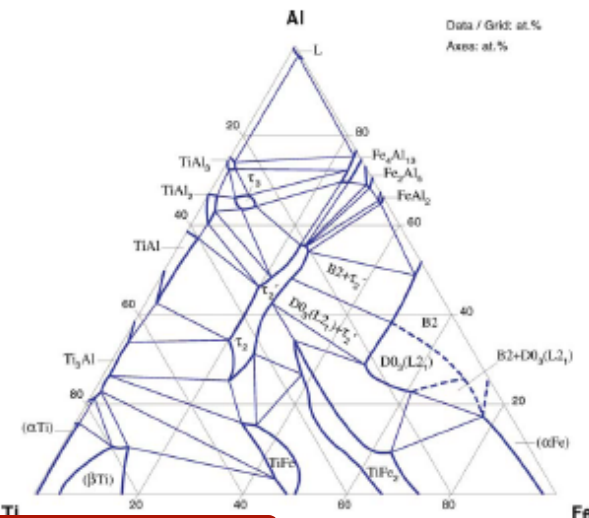


Fig. 14. Al-Fe-Ti Isothermal section at 800°C

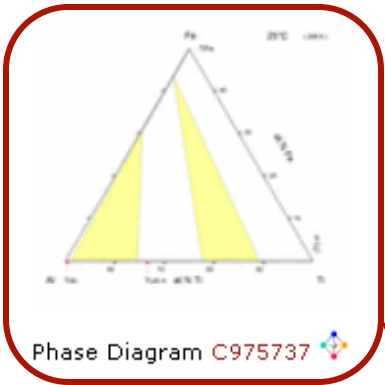
**„Landolt-Börnstein“ Content:  
Search Hit Highlighted**

**Al-Fe-Ti, ternary phase diagram, isothermal section**

Element System: Al-Fe-Ti

**Inorganic Solid Phases · phase diagrams**

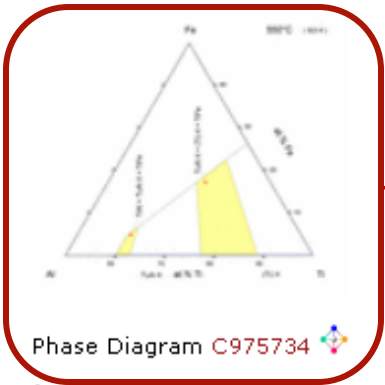
Diagram details: Al conc.[0-50 at.%] vs. Fe conc.[0-50 at.%] vs. Ti conc.[50-100 at.%]



Phase Diagram C975737

from:  
Seibold A., Seibold A., Seibold A., Seibold A.: "Phasengleichgewichte in den ternären Systemen Ti-Fe-O und Ti-Al-Fe", Z. Metallkd. **72** (1981) 712-719.

Diagram details: Al conc.[0-30 wt.%] vs. Fe conc.[0-30 wt.%] vs. Ti conc.[70-100 wt.%]



Phase Diagram C975734

from:  
Volkova M.A., Kornilov I.I.: "PHASE EQUILIBRIUM IN TITANIUM-RICH Ti-Al-Fe ALLOYS", Russ. Metall. **3** (1970) 134-137.

**„Inorganic Solid Phases“:  
Overview of Relevant  
Data Sheets**

Diagram details: Al conc.[0-45 at.%] vs. Fe conc.[0-27 at.%] vs. Ti conc.[55-100 at.%]

Element System: Al-Fe-Ti

Ternary Phase Diagram

Investigated Concentration Ranges: Al conc.[0-50 at.%) vs. Fe conc.[0-50 at.%) vs. Ti conc.[50-100 at.%)

**Constitution Entry Summary**

<b>System</b>	Al-Fe-Ti
<b>Diagram Type</b>	ternary, isothermal section
<b>Concentration Range</b>	partial composition; Ti-TiAl-TiFe / 50-100 at.% Ti
<b>Temperature</b>	298 K
<b>Investigation</b>	experimental
<b>APDIC</b>	non-APDIC

**Reference**

Seibold A.: "Phasengleichgewichte in den ternären Systemen Ti-Fe-O und Ti-Al-Fe", Z. Metallkd. 72 (1981) 712-719.

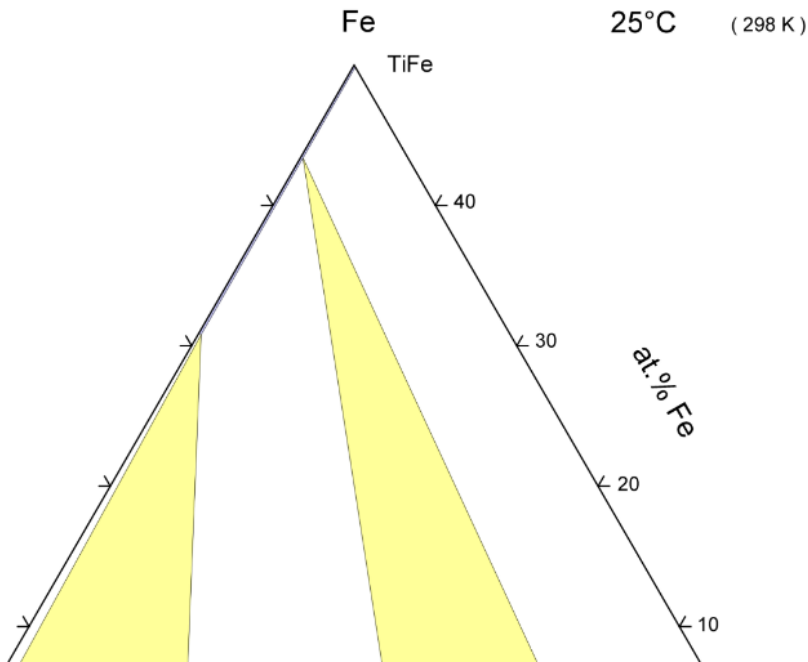
**Phase Diagram**

**Published Phase Diagram**

<b>Figure</b>	19
<b>Title</b>	Al-Fe-Ti isothermal section at 25 °C
<b>Parameters</b>	Al conc.[0-50 at.%) vs. Fe conc.[0-50 at.%) vs. Ti conc.[50-100 at.%)

**Redrawn Phase Diagram**

<b>Concentration range</b>	partial composition; Ti-TiAl-TiFe
<b>Enlargement</b>	0.6



**Actual Data Sheet**

Results 1 - 8 of 8 Documents




Thermophysical Properties

**Methanol / Water** ↻

Metadata - Substance: Methanol ... Water ...

Molecules and Radicals > Radicals, Reaction Kinetics > Oxyl-, Alkoxy-, Carbonyloxy-, Phenoxy- and Peroxy-Radicals in Liquids > Peroxy and related radicals > Alkylperoxy radicals > Radical-molecule reactions

**Electron transfer reactions with organic compounds** 📄 i

**Fulltext:** (10%) (30%) (60%) (80%) (95%) **methanol water** + acetonitrile (84%) water + ...

Molecules and Radicals > Radicals, Magnetism > Free Organic C- and N-Centered and Nitroxide Radicals > Nitroxide Radicals and Polynitroxides (Update) > Cyclic nitroxides > Pyrrolidiny-N-oxyls

**Trisubstituted pyrrolidiny-N-oxyls** 📄 i

**Fulltext:** oxidation ESR / 298 Toluene **Methanol Water** Photolysis of oxygen-ated ... and benzo-phenone ESR / 298 **Methanol Water** Photolysis of glycine, ...

Molecules and Radicals > Radicals, Magnetism > Free Organic C- and N-Centered and Nitroxide Radicals > Nitroxide Radicals and Polynitroxides (Update) > Cyclic nitroxides > Pyrrolidiny-N-oxyls

**Tetrasubstituted pyrrolidine-N-oxyls** 📄 i

**Fulltext:** ESR / 298 Benzene **Methanol Water**, pH 1.0 (pH 14.0) Reaction of ...

Molecules and Radicals > Radicals, Reaction Kinetics > Proton and Electron Transfer, Biradicals and Excited Radicals in Liquids > Biradicals > Unimolecular biradical processes

**Reactions of 1,4-biradicals to yield molecular products**

**Fulltext:** acetonitrile THF 2-propanol **methanol water** dimethylsulfo

Advanced Materials > Proteins > Structural and Physical Data of proteins and related substances

**Tables 1...9** 📄 i

**Fulltext:** 75 % dioxane water 1 M NaCl **methanol water** water chloroform isopropyl ...

Molecules and Radicals > Radicals, Magnetism > Free Organic Anion and Cation Radicals, Polyradicals > Anion Radicals > Semiquinones and related species > Semiquinone anion radicals > Naphthosemiquinones

**“Thermophysical Properties”:  
Data Sheet Found**

**Methanol / Water**

**Thermophysical Data in the Dortmund Data Bank**

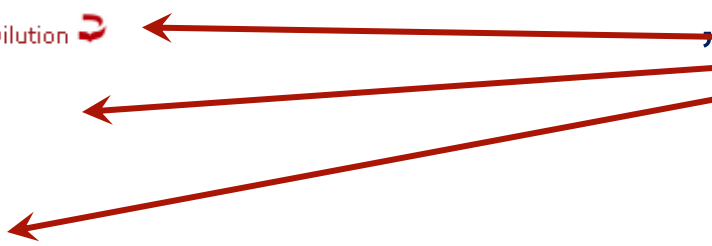
**Components**

No.	Formula	Molar Mass	CAS Registry Number	Name
1	CH <sub>4</sub> O	32.042	67-56-1	Methanol
2	H <sub>2</sub> O	18.015	7732-18-5	Water

**List of Available Properties**

- Activity Coefficients at Infinite Dilution ↻
- Azeotropic Data ↻
- Excess Heat Capacities ↻
- Heat Capacities ↻
- Excess Enthalpies ↻
- Solid-Liquid Equilibria ↻
- Excess Volumes ↻
- Densities ↻
- Vapor-Liquid Equilibria ↻

**Thermophysical Properties“: Overview of Available Data**



Dortmund Data Bank Edition - 2009. Integrated in SpringerMaterials - 2010  
 Dortmund Data Bank by DDBST - Dortmund Data Bank Software and Separation Technology GmbH  
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**Methanol / Water**  
**Activity Coefficients at Infinite Dilution**
**Components**

No.	Formula	Molar Mass	CAS Registry Number	Name
1	CH <sub>4</sub> O	32.042	67-56-1	Methanol
2	H <sub>2</sub> O	18.015	7732-18-5	Water

**Water solved in Methanol**

Temperature [K]	Activity Coeff. at Inf. Dil.	Measurement Method	Source
298.15	1.6300	CPEQ	4
298.15	4.1800	OTHR	2
307.75	1.6000	EBUL	1
313.15	2.2000	STAT	3
317.85	1.8000	EBUL	1
323.15	3.2600	OTHR	2
327.85	1.7000	EBUL	1
333.15	1.7700	CPEQ	4
337.05	1.7600	EBUL	1
373.15	1.8700	CPEQ	4

[Download Data Table as CSV File](#)
**List of measurement methods**

Short Term	Explanation
CPEQ	calculated from phase equilibria
EBUL	ebullimetry
OTHR	other techniques
STAT	static method

**List of References**

1	Bergmann D.L., Eckert C.A.: Measurement of Limiting Activity Coefficients for Aqueous Systems by Differential Ebullimetry. Fluid Phase Equilib. 63 (1991) 141-150
2	Schmidt T.W.: Determination of Infinite Dilution Activity Coefficients (gamma infinite) using Molecular Beams. US-Patent (1980) 1-16
3	Belousov V.P., Makarova N.I., Sabylin I.I.: Thermodynamic Properties of the Water-Ethanol-Dioxan Ternary System at 25°C. Viniti (1973) 1-10
4	Pierotti G.J., Deal C.H., Derr E.L.: (Addendum included). Ind.Eng.Chem. 51 (1959) 95-102

**Methanol solved in Water**

Temperature [K]	Activity Coeff. at Inf. Dil.	Measurement Method	Source
273.35	1.2300	STAT	26
283.15	1.4000	STAT	26
283.15	1.4100	DILU	22
293.15	1.5500	DILU	22
293.15	1.5500	STAT	26
293.15	1.6800	STAT	16
293.15	2.6900	GLCI	1
297.45	2.1200	GLCI	2
298.15	1.5300	HENR	15
298.15	1.5300	CPEQ	23
298.15	1.5800	GLCI	17
298.15	1.6400	GLCR	8
298.15	1.6400	DILU	6

**Actual Data Sheet**
**Data available as CSV File**

Go Advanced Search

Home Bookshelf Periodic Table Help For Librarians Feedback

- Al-Fe-Ti
- Al-Fe-Ni-Ti
- Al-Fe-O-Ti
- Al-Fe-Ti-V
- Al-B-Fe-Ni-Ti
- Al-Cr-Fe-O-Ti
- Al-Fe-Mg-O-Ti
- Al-Fe-Na-O-Ti
- Al-Fe-O-Si-Ti
- Al-B-Fe-Mg-O-Ti
- Al-Ca-Fe-Mg-Mn-Ti
- Al-Ca-Fe-O-Si-Ti
- Al-Ca-Fe-O-Ti-Zr
- Al-Ce-Fe-O-Si-Ti
- Al-Co-Cu-Fe-Ni-Ti
- Al-Cr-Fe-Mg-O-Ti
- Al-Cr-Fe-Nb-Ni-Ti
- Al-Fe-H-O-Si-Ti
- Al-Fe-Mg-O-Si-Ti
- Al-Fe-Mn-O-Pb-Ti
- Al-Fe-Mo-O-Sr-Ti
- Al-Fe-Nd-O-Si-Ti
- Al-Fe-O-Pr-Si-Ti
- Al-Ba-Fe-Mn-Ni-O-Ti
- Al-Ca-Ce-Fe-O-Si-Ti
- Al-Ca-Cr-Fe-O-Si-Ti
- Al-Ca-Fe-H-O-Si-Ti
- Al-Ca-Fe-La-O-Si-Ti

## Search for Element Systems

Select elements by clicking on the symbols. Deselect elements by clicking a second time.

**Your Selection**  
Al-Fe-Ti

	1 IA	2 IIA	3 IIIB	4 IVB	5 VB	6 VIB	7 VIIB	8 VIII	9 VIII	10 VIII	11 IB	12 IIB	13 IIIA	14 IVA	15 VA	16 VIA	17 VIIA	18 VIIIA			
1	H	He																	He	K	
2	Li	Be											B	C	N	O	F	Ne		L	
3	Na	Mg											Al	Si	P	S	Cl	Ar		M	
4	K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	Se	Br	Kr		N	
5	Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Ag	Cd	In	Sn	Sb	Te	I	Xe		O	
6	Cs	Ba	*	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At	Rn		P	
7	Fr	Ra	**	Rf	Db	Sg	Bh	Hs	Mt	Ds	Rg	Cn									Q
*		La	Ce	Pr	Nd	Pm	Sm	Eu	Gd	Tb	Dy	Ho	Er	Tm	Yb	Lu					
**		Ac	Th	Pa	U	Np	Pu	Am	Cm	Bk	Cf	Es	Fm	Md	No	Lr					

Available Content



Substances, Properties, ... | Bibliographic References | Help | Close

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- Advanced
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- Solid Phase
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- Chemical

urea
urea (57-13-6; CH <sub>4</sub> N <sub>2</sub> O)
Urea-d4 (1433-11-0; CD <sub>4</sub> N <sub>2</sub> O)
Urea-d4 (1433-11-0; CD <sub>4</sub> N <sub>2</sub> O)
Urea-12C (57-13-6; CH <sub>4</sub> N <sub>2</sub> O)
Urea-phil (57-13-6; CH <sub>4</sub> N <sub>2</sub> O)
(2H <sub>4</sub> )urea (1433-11-0; CD <sub>4</sub> N <sub>2</sub> O)
Ethyl-urea (625-52-5; C <sub>3</sub> H <sub>8</sub> N <sub>2</sub> O)
butyl-urea (592-31-4; C <sub>5</sub> H <sub>12</sub> N <sub>2</sub> O)
butyl-urea (592-31-4; C <sub>5</sub> H <sub>12</sub> N <sub>2</sub> O)
Acetyl-urea (591-07-1; C <sub>3</sub> H <sub>6</sub> N <sub>2</sub> O <sub>2</sub> )
Benzyl-urea (538-32-9; C <sub>8</sub> H <sub>10</sub> N <sub>2</sub> O)
Phenyl-urea (64-10-8; C <sub>7</sub> H <sub>8</sub> N <sub>2</sub> O)
Urea, thio- (62-56-6; CH <sub>4</sub> N <sub>2</sub> S)
methyl-urea (598-50-5; C <sub>2</sub> H <sub>6</sub> N <sub>2</sub> O)
1-Ethyl urea (625-52-5; C <sub>3</sub> H <sub>8</sub> N <sub>2</sub> O)
Benzoyl urea (614-22-2; C <sub>8</sub> H <sub>8</sub> N <sub>2</sub> O <sub>2</sub> )

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Substances / Molecular Formulas / Element Systems / CAS Registry Numbers

11

- 110-00-9 (furan)
- 110-01-0 (tetrahydrothiophene)
- 110-02-1 (thiophene)
- 110-05-4 (di-*tert*-butyl peroxide)
- 110-06-5 (di-*tert*-butyl disulfide)
- 110-12-3 (5-methyl-hexan-2-one)
- 110-13-4 (hexane-2,5-dione)
- 110-14-5 (Succinamid)
- 110-15-6 (butanedioic acid)
- 110-16-7 (*cis*-butenedioic acid)
- 110-17-8 (Fumaric acid)
- 110-18-9 (*N,N,N',N'*-Tetramethyl-ethane-1,2-diamine)
- 110-19-0 (acetic acid isobutyl ester)
- 110-20-3 (Hydrazinecarboxamide, 2-(1-methylethylidene)-)
- 110-21-4 (Diharnstoff)
- 110-22-5 (Peroxide, diacetyl)

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- Chemical Safety**

## Search in

# REACH, GHS, RoHS, WEEE

Substances / Molecular Formulas / Element Systems / CAS Registry Numbers

REACH - **R**egistration, **E**valuation, **A**uthorization and **R**estriction of **C**hemicals  
GHS - **G**lobally **H**armonized **S**ystem  
RoHS - **R**estriction of **H**azardous **S**ubstances  
WEEE - **W**aste from **E**lectrical and **E**lectronic

- Particles, Nuclei and Atoms
- Molecules and Radicals
- Electronic Structure and Transport
- Magnetism
- Semiconductivity
- Superconductivity
- Crystallography
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**Search in**  
**REACH, GHS, RoHS, WEEE**

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- benzene (C<sub>6</sub>H<sub>6</sub>)
- benzene-d<sub>6</sub> (C<sub>6</sub>D<sub>6</sub>)
- Benzene-1,4-d<sub>2</sub> (C<sub>6</sub>H<sub>4</sub>D<sub>2</sub>)
- benzenemethanol (C<sub>7</sub>H<sub>8</sub>O)
- Cyanato-benzene (C<sub>7</sub>H<sub>5</sub>NO)
- benzene selenol (C<sub>6</sub>H<sub>6</sub>Se)
- butoxy-benzene (C<sub>10</sub>H<sub>14</sub>O)
- dimethyl-benzene (C<sub>8</sub>H<sub>10</sub>)
- propoxy-benzene (C<sub>9</sub>H<sub>12</sub>O)
- Benzene, hexaiodo- (C<sub>6</sub>I<sub>6</sub>)


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European regulations regarding benzenemethanol (C<sub>7</sub>H<sub>8</sub>O)


**Name** benzenemethanol **Formula:** C<sub>7</sub>H<sub>8</sub>O  
**CAS-RN** 100-51-6 **Molecular Weight:** 108.138 g/mol  
**EG-Index:** 603-057-00-5 (93/72/EEC)  
**EINECS:** 202-859-9 (EINECS)

## Hazard Information (Dangerous Substances Directive 67/548/EEC)

<b>Hazard symbol</b>		93/72/EEC
<b>R-Phrase</b>	Xn Harmful 20/22 R20/22 Harmful by inhalation and if swallowed.	93/72/EEC
<b>S-Phrase</b>	(2-)26 S2 Keep out of the reach of children. S26 In case of contact with eyes, rinse immediately with plenty of water and seek medical advice.	93/72/EEC

## GHS classification (Globally Harmonized System)

Regulation on Classification, Labelling and Packaging of Substances and Mixtures (CLP)

<b>Signal Word</b>	Warning	EC/1272/2008
<b>Pictogram</b>		
<b>Hazard Statements</b>	H332 Harmful if inhaled. H302 Harmful if swallowed.	
<b>Precautionary Statements</b>	P261 Avoid breathing dust/fume/gas/mist/vapours/spray. P264 Wash ... thoroughly after handling. P270 Do not eat, drink or smoke when using this product. P271 Use only outdoors or in a well-ventilated area. P301+P312 IF SWALLOWED: Call a POISON CENTER or doctor/physician if you feel unwell. P304+P340 IF INHALED: Remove victim to fresh air and keep at rest in a position comfortable for breathing. P312 Call a POISON CENTER or doctor/physician if you feel unwell. P330 Rinse mouth. P501 Dispose of contents/container to ...	
<b>Classification</b>	Acute Tox. 4 * Acute Tox. 4 *	
<b>EC-Reference</b>	603-057-00-5	

Actual Safety Document

## European Chemicals Agency (ECHA) pre-registration

<b>Pre-registered substance</b>	EC-Number: 202-859-9 CAS-Number: 100-51-6 Name: benzyl alcohol Synonym: - Registration Date: 30/11/2010	ECHA pre
---------------------------------	---	----------

## References

**93/72/EEC**  
**Short:** 93/72/EEC  
**Title:** Commission Directive 93/72/EEC of 1 September 1993 adapting to technical progress for the nineteenth time Council Directive 67/548/EEC on the approximation of the laws, regulations and administrative provisions relating to the classification, packaging and labelling of dangerous substances  
**Author:** Council of the European Economic Community  
**Source:** *Official Journal of the European Communities*  
**Volume:** L 258 A (16.10.1993)  
**Page:** 1-1409  
**Year:** 1993  
**Keyword:** hazardous materials; hazard classification  
**Internet Resource:** <http://eur-lex.europa.eu/LexUriServ/LexUriServ.do?uri=CELEX:31993L0072:EN:HTML>  
**Publish Date:** 1993/10/16  
**Short:** EC/1272/2008  
**Title:** REGULATION (EC) No 1272/2008 OF THE EUROPEAN PARLIAMENT AND OF THE COUNCIL of 16 December 2008 on classification, labelling and packaging of substances and mixtures, amending and repealing Directives 67/548/EEC and 1999/45/EC, and amending Regulation (EC) No 1907/2006  
**Author:** European Parliament; Council of the European Union  
**Source:** *Official Journal of the European Union*

EC/1272/2008

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Al-Fe-Ti "yield strength"

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Improvement in mechanical properties of NiAl matrix composites ...

by YX Lu - 1996 - Related articles
With Ni, Al, Fe, Ti and C elemental powders, the NiAl(Fe) matrix composites have .... The yield strength increases with increasing volume fraction of TiC ...
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Microstructure and compression properties of TiCp/NiAl(Fe ...

by YX Lu - 1997 - Cited by 2 - Related articles
It shows that the yield strength (0.2 YS) decreases with increasing the ... Conclusions (1)
With Ni, Al, Fe, Ti and C elemental powders, NiAl(Fe) matrix ...
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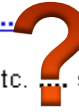
ATOMISTIC STUDIES OF DEFORMATION MECHANISM OF NANOCRYSTALLINE AL ...

by T Uesugi - 2007
It is to be noted that the yield strength of the Al-Ti solid solution is ... atomic solute cluster, contributes to the high yield strength of Al-Fe alloys. ...
www.scientific.net/MSF.561-565.977.pdf



lower steel tapping temperature at LD/BOF which impacts ...

File Format: PDF/Adobe Acrobat - Quick View
to requirements of steel properties e.g. yield strength, ductility, formability etc. .... stringent quality requirement and Al in Fe-Ti can give rise to ...
www.pyrometallurgy.co.za/InfaconXI/082.pdf



Development of Low Cost Powder Metallurgy Process of Titanium ...

File Format: PDF/Adobe Acrobat - Quick View
High fatigue strength with respect to yield strength can be cited as one characteristic of the Ti-Fe-Al ternary sintered alloys. Fig. 5 ...
www.nsc.co.jp/en/tech/report/pdf/8515.pdf - Similar



Synthesis and Characterization of Mechanically Alloyed and HIP ...

by SM Zhu - Related articles
The weakening of Fe, Ti and Al peaks indicates that dissolution due to ... yield strength usually exhibited by cast Fe3Al-based alloys is not observed. ...
www.mrs.org/s\_mrs/bin.asp?CID=2398&DID=56263&DOC=FILE...



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Sort by: **Relevance** ▾ Date Title

Did you mean '**al feti yield strength**' rather than '**Al-Fe-Ti yield strength**'?

- Journal Article

**Formation of simple crystal structures in Cu-Co-Ni-Cr-Al-Fe-Ti-V alloys with multiprincipal metallic elements**

Jien-Wei Yeh, Su-Jien Lin, Tsung-Shune Chin, Jon-Yaw Gan and Swe-Kai Chen, *et al.*  
*Metallurgical and Materials Transactions A*, 2004, Volume 35, Number 8, Pages 2533-2536

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- Journal Article

**Petrology, chronology and isotope geochemistry of the proterozoic amphibolites from Xiangshan, central Jiangxi province, China**

Hu Gongren, Zhang Bangtong and Yu Ruilian  
*Chinese Journal of Geochemistry*, 1999, Volume 18, Number 2, Pages 139-149

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- Book Chapter

**Nanoparticles and Polyelectrolytes**

*Springer Laboratory*, 2007, *Polyelectrolytes and Nanoparticles*, Pages 47-71

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Refine

Multiphase Systems > Ternary Alloys > Phase Diagrams, Crystallography and Thermodynamics > Light Metal Systems > Aluminum (Al-X-Y) Ternary Alloys

Al-Fe-Ti

**Metadata - Substance: Al-Fe-Ti ... Metadata - Element System: Al-Fe-Ti ... Fulltext:** Equilibrium Diagram of the **Al-Fe-Ti** System and the Segregation of ... electrical resistivity and **yield strength** measurements [1940Fin] and by ... of Al promote an increasing **strength** and creep resistance when ... Design of the High **Strength**, Creep-Resistant Aluminium ...

Multiphase Systems > Ternary Alloys > Phase Diagrams, Crystallography and Thermodynamics > Iron Systems > Selected Systems from Al-B-Fe to C-Co-Fe

Aluminium - Iron - Titanium

**Metadata - Substance: Al-Fe-Ti ... Metadata - Element System: Al-Fe-Ti ... Fulltext:** 1. Investigations of the **Al-Fe-Ti** Phase Relations, Structures ... Temperature dependence of **yield stress** [1991Kum] Mechanical ... Temperature dependence of **yield stress** [1991Wu] Mechanical ... tests Tensile and fracture **strengths** [2003Ste] Mechanical tests ...

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# Use Cases

Aircraft engineer needs new alloy for construction of stronger winglets: e.g., Al-Fe-Ti



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Your Selection  
Al-Fe-Ti

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IA	IIA	IIIB	IVB	VB	VIB	VIIb	VIIIb	VIIIb	VIIIb	IB	IIB	IIIA	IVA	VA	VIA	VIIA	VIIIA	
1	H	He															He	
2	Li	Be										B	C	N	O	F	Ne	
3	Na	Mg										Al	Si	P	S	Cl	Ar	
4	K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	Se	Br	Kr
5	Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Au	Hg	Tl	Pb	Bi	Po	At	Rn
6	Cs	Ba	**	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At	Rn
7	Fr	Ra	**	Rf	Db	Sg	Bh	Hs	Mt	Ds	Rg	Cn						
*	La	Ce	Pr	Nd	Pm	Sm	Eu	Gd	Tb	Dy	Ho	Er	Tm	Yb	Lu			
**	Ac	Th	Pa	U	Np	Pu	Am	Cm	Bk	Cf	Es	Fm	Md	No	Lr			

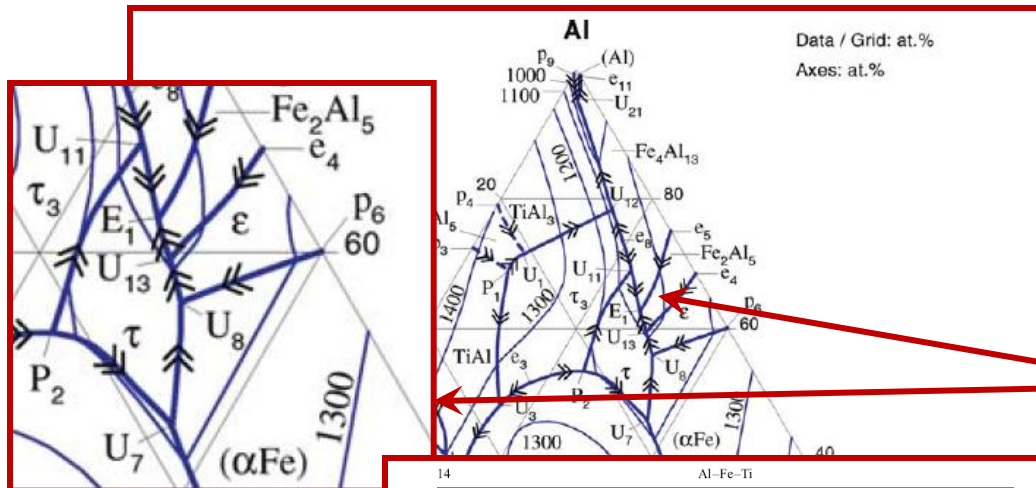
1. Selects components by clicking on elements in periodic table
2. Available content shown

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## Typical Usage Situation



1. Opens phase diagram
2. Selects desired material from diagram
3. Collects desired materials properties

Al-Fe-Ti

Phase/ Temperature Range [°C]	Pearson Symbol/ Space Group/ Prototype	Lattice Parameters [pm]	Comments/References
		$a = 496.61$ $c = 806.28$	[2006Yan], $Ti_{33.3}Fe_{33.3}Al_{33.4}$ annealed at 1000°C
		$a = 503.66$ $c = 819.71$	[2006Yan], as-cast $Ti_{34}Fe_{17}Al_{49}$
TiFe ≤ 1317	$cP2$ $Pm\bar{3}m$ CsCl	$a = 297.6$	solid solubility ranges from 49.8 to 51.8 at.% Ti [V-C]
* $\tau_1$ , TiFe <sub>2</sub> Al	$cF16$ $Fm\bar{3}m$ Cu <sub>2</sub> AlMn	$a = 587.9$	[1983Bus], annealed at 900°C for 14 days
* $\tau_2$	$cF*$ $F\bar{4}3m$	$a = 1211.0$ $a = 1209.59$	[1995Pal], at $Ti_{50.9}Fe_{24.5}Al_{24.6}$ [2006Gry], $Ti_{13}Fe_{24}Al_{33}$ annealed at 900°C; both X-ray and neutron diffraction data
* $\tau_2'$	$cF116$ $Fm\bar{3}m$ Th <sub>6</sub> Mn <sub>23</sub>	$a = 1199.0$ $a = 1182.0$ $a = 1203.8$ $a = 1207.6$ $a = 1209.9$ $a = 1211.0$ $a = 1189.0$	[1967Mar, 2000Mab] [1981Sei] [1995Pal], at $Ti_{30.8}Fe_{21.4}Al_{47.8}$ [1999Gor], at $Ti_{38.4}Fe_{23}Al_{38.6}$ [1999Gor], at $Ti_{42.2}Fe_{23.1}Al_{34.7}$ [1999Lev] [2003Gry] at $Ti_{20.3}Fe_{23.7}Al_{56}$
	filled Th <sub>6</sub> Mn <sub>23</sub>	$a = 1209.2$	[2003Gry] at $Ti_{42}Fe_{23.3}Al_{34.7}$
	filled Th <sub>6</sub> Mn <sub>23</sub>	$a = 1199.44$	[2006Gry], $Ti_{22}Fe_{23}Al_{35}$ annealed at 900°C; neutron diffraction data

# Chemical reactor burst due to high vapor pressure



# SpringerMaterials Typical Usage Situation

40 2.2 Hydrocarbons, C<sub>6</sub> to C<sub>9</sub> [Ref. p. 261]

Phase	A Antoine constants A, (°C) B [K], (K) C [K], (°C)	T <sub>c</sub> [K]	Range [K], Rating	T <sub>c</sub> [K], P <sub>c</sub> [kPa]	Ref.
387	C <sub>8</sub> H <sub>18</sub>	Octane			111-65-9
l-g	6.56398 1606.62 -42.89	217.204	216.4/204 B	398.83/101.325	90-90-90
l-g	6.05075 1356.36 -63.515	288.423	284.407 A	284.407 A	90-90-90
l-g	6.05075 1356.36 -63.515	423.568	407/568.4 B	407/568.4 B	90-90-90
	(2.88414) (948.2)	(-73346)			
388	C <sub>8</sub> H <sub>18</sub>	2,2,3,3-Tetramethylbutane			694-82-1
cr-g	5.91839 1652.6 -46.09	252.372	242/374 B	379.44/101.325	90-90-90
l-g	5.9042 1270.1 -53.65	372.406	374/416 B	374/416 B	90-90-90
389	C <sub>8</sub> H <sub>18</sub>	2,2,3-Trimethylpentane			564-02-3
l-g	5.94826 1293.84 -54.795	284.408	270/400 A	382.99/101.325	90-90-90
l-g	5.94826 1293.84 -54.795	408.563	400/563.5 B	400/563.5 B	90-90-90
	(2.45345) (162.4)	(-65383)			
390	C <sub>8</sub> H <sub>18</sub>	2,2,4-Trimethylpentane			549-84-1
l-g	6.35751 1447.78 -36.53	180.272	180/272 B	372.90/101.325	90-90-90
l-g	5.93646 1257.85 -52.383	272.980	272/400 A	272/400 A	90-90-90
l-g	5.93646 1257.85 -52.383	398.553	400/543.9 B	400/543.9 B	90-90-90
	(1.12261) (134.5)	(12998)			
391	C <sub>8</sub> H <sub>18</sub>	2,3,3-Trimethylpentane			560-21-4
l-g	5.96421 1325.81 -52.989	287.408	277/403 A	387.92/101.325	90-90-90
l-g	5.96421 1325.81 -52.989	408.573	402/573.5 B	402/573.5 B	90-90-90
	(2.3795) (76.3)	(1851)			
392	C <sub>8</sub> H <sub>18</sub>	2,3,4-Trimethylpentane			565-75-3
l-g	6.35762 1507.04 -38.35	216.298	205/298 B	386.62/101.325	90-90-90
l-g	5.977 1314.31 -55.669	288.408	288/400 A	288/400 A	90-90-90
l-g	5.977 1314.31 -55.669	408.566	400/566.4 B	400/566.4 B	90-90-90
	(2.38574) (169.4)	(-44867)			
393	C <sub>9</sub> H <sub>12</sub>	Indane, (1H-indene)			95-13-6
l-g	6.34410 1749.215 -52.375	297.457	290/480 B	455.57/101.325	90-90-90
					6-Chance
394	C <sub>9</sub> H <sub>12</sub>	Indan			496-11-7
l-g	6.11230 1577.321 -66.828	375.465	370/470 A	450.92/101.325	90-90-90
395	C <sub>9</sub> H <sub>12</sub>	Isopropylbenzene			95-83-0
l-g	7.13460 2234.172 -2.336	395.438	290/440 C	437.34/101.325	90-90-90
					90-90-90
396	C <sub>9</sub> H <sub>12</sub>	2-Methylstyrene			611-15-4
l-g	6.27022 1624.066 -62.132	305.985	300/390 B	370.20/10	90-90-90
397	C <sub>9</sub> H <sub>12</sub>	3-Methylstyrene			109-90-1
l-g	6.36538 1682.941 -56.908	314.442	314/330 C, 330/445 B	442.93/101.325	90-90-90

Ref. p. 12 | 1 Introduction | 7

### 1.3.2 Empirical Vapor Pressure Equations

During the past century many empirical mathematical functions have been used to relate vapor pressure to temperature; most are modifications of Eq. (1.7). These functions have several parameters that are characteristic of the compound. Curve fits off experimental data, usually by minimizing the sum of the squares of the deviations between the calculated and observed pressures or temperatures (least squares criterion), provide these parameters. The first and most widely used of these equations is the Antoine equation [168-ant, 16-6to]. The original form is,

$$\log P = A - B / (C + T) \quad (1.8)$$

Sometimes the natural logarithm is used instead of the base-10 logarithm or Celsius temperature is used instead of Kelvin. When  $C = 0$  (for  $T$  in kelvins) Eq. (1.8) is identical to Eq. (1.7). The *Thermodynamics Research Center Thermodynamic Tables - Hydrocarbons [20-4to]* and *Nonhydrocarbons [20-4to]* - use an extended version of the Antoine equation:

$$\log P = A - B / (C + T)^1 + 0.434294 T^2 + E T^3 + F T^4 \quad (1.9)$$

where  $n$ ,  $E$ , and  $F$  are additional adjustable parameters.  $T_c$  is the critical temperature,  $T_0$  is the lower boundary temperature and  $\chi = (T - T_0)/T_c$ . Examples of functions obtained by adding terms to Eq. (1.7) are the polynomial in temperature used in the *International Critical Tables [26-ant]*,

$$\ln P = A + BT^{-1} - CT + DT^2, \quad (1.10)$$

the Chebyshev polynomial [70-ambcoo]

$$T \ln P = a_0 + \sum_{i=1}^n a_i E_i(\chi) \quad (1.11)$$

in which  $E_i(\chi)$  is a Chebyshev polynomial in  $\chi$  of degree  $i$  (the advantage of this is that the  $E_i$  functions are orthogonal), the Kirchoff-Ramaline equation [48-4to],

$$\chi = [2T - (T_{\text{sat}} - T_{\text{sat}})] / (T_{\text{sat}} - T_{\text{sat}}) \quad (1.12)$$

and the Frost-Kalkwarf equation [53-froka]

$$\ln P = A + BT^{-1} + C \ln T, \quad (1.13)$$

(same form as Eq. (1.6)), the Plank-Riedel equation [48-plank]

$$\ln P = A + BT^{-1} + C \ln T + DP^e, \quad (1.14)$$

and the Frost-Kalkwarf equation [53-froka]

$$\ln P = A + BT^{-1} + C \ln T - DPT^{-2} \quad (1.15)$$

Another popular type of function is the Cox equation [36-cox]:

$$\ln (PP_c^*) = A (1 - T/T_c) \quad (1.16)$$

where  $A$  is a function of temperature often chosen to be

$$\ln A = a + bT + cT^2 \quad (1.17)$$

Wagner and others [76-wag, 73-wag-1, 77-wag, and 86-amb-1] have proposed a series of related equations. The simplest is

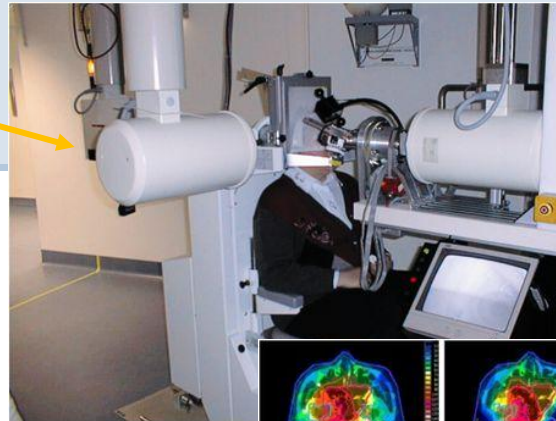
$$\ln (PP_c^*) = (A - Bt^{1.5} - Ct^3 - Dt^6) / T, \quad (1.18)$$

where  $t = 1 - T/T_c$ ,  $P_c$  is the critical pressure and  $T_c$  is the critical temperature. One of the variations [76-wagwe] is:

$$\ln (PP_c^*) = (A - Bt^{1.5} - Ct^3 - Dt^6 - Et^9) / T, \quad (1.19)$$

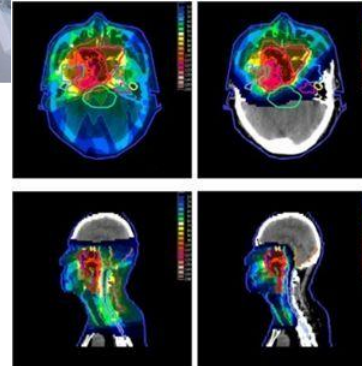
1. Read fundamentals
2. Select desired vapor pressure from table

Manufacturer of medical equipment designs a new apparatus for tumor therapy



# SpringerMaterials

## Typical Usage Situation



Ref. [77] 1.7 Production of atomic defects by irradiation Figs. p. 37

Table 5 (continued)

Metal	$\phi_d$ [ $\mu\text{hr/s/a.c.}$ ]	Method	Ref.	Metal	$\phi_d$ [ $\mu\text{hr/s/a.c.}$ ]	Method	Ref.
Mg	$\geq 0.8$	ERM	69 F	Tb	135–20	ERM	80 D 1
	4–5	FRM	72 O	Tb	15	FRM	75 G 2
	2 <sup>1)</sup>	FRM	75 L 1	(19 <sup>1)</sup> )	FRM	75 L 1	
Mo	2	HS-LPM/ERM	52 E 1	Ti	18	ERM	72 S 1
	4–5	ERM	62 L 1	(10 <sup>1)</sup> )	ERM	75 L 1	
	10	FRM	75 R	Tm	140–50	FRM	80 D 1
	13–2	ERM-S	75 M 5	U	22	ERM-LPM	87 W
Nb	12–4	HS-LPM/ERM	78 e	V	5.2–7.5	ERM	75 C 1
	16	FRM	75 I 1		21.6 <sup>2)</sup>	ERM	80 7
	16–23	HS-LPM/ERM	78 b		22–7 <sup>1)</sup>	FRM	85 D 3
Ni	135–135	ERM	85 D 1	W	7.5–16.0	ERM-S	78 M
Ni	6	ERM	75 L 1		14	FRM	75 L 1
	7.1±0.8	HS-LPM/FRM	85 B		28	FRM	80 K
Yb	$\leq 1$	ERM	74 B 2		27–6 <sup>1)</sup>	ERM	85 D 3
	20	ERM	75 L 1	V	30–20	ERM	80 D 2
Pd	2±1	ERM	67 1	Yb	75–25	ERM	80 D 2
	9.7±0.5 <sup>1)</sup>	FRM	85 D 2	Zn	5	FRM	71 M 2
Pr	135+35	ERM	85 D 1		4.2+5	ERM	73 M 4
Pt	9.5±0.5	ERM-S	73 J 2		20–3	ERM-S	73 M 2
	20	ERM	74 V 1		15–5	ERM-S	77 v
Se	69	FRM	85 D 4		12–5	HS-LPM/ERM	78 E
Sm	140 130	ERM	80 D 2		15.3	HS-LPM/ERM	79 E 2
Sn(2)	1.1±0.2	ERM	75 M 1	Zr	35.0	ERM	75 N
	4+2 <sup>2)</sup>	ERM	85 D 2		40	ERM	71 B 2
Ta	17±3	ERM-S	72 1		35–8	HS-LPM/ERM	82 F 1
	16±3	ERM-S	79 B 3				

<sup>1)</sup> Reevaluation of preceding data  
<sup>2)</sup> Estimated value.

### 1.7.1.2 Ion irradiation

For nonrelativistic particles the maximum transferred energy is:

$$T_{max} = E \cdot f$$

$$w_{ph} = \gamma - 4M_1^2 M_2^2 / (M_1^2 + M_2^2)^2 \quad (2)$$

The differential cross section for nonrelativistic charged particles is given by the Rutherford cross section:

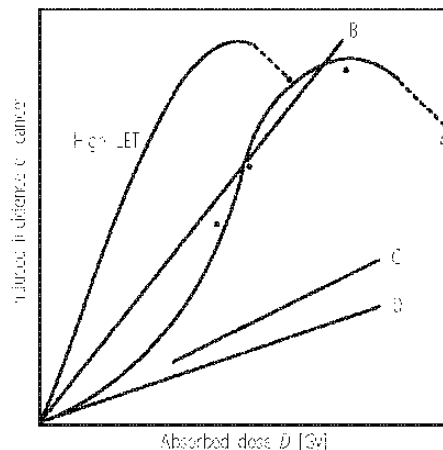
$$d\sigma/dT = \pi M_2^2 / (M_1^2 (2Z_1 Z_2 e^2)^2 (1-f)^2) \quad (3)$$

Corrections to the Rutherford cross section must be applied at low energy transfers to account for shielding of the nucleus charge by atomic electrons [68.1]. An analytical expression for this corrected cross section was derived in [72 W 1]. At high energies modifications of the Rutherford cross section are necessary, when inelastic nuclear scattering and nuclear reactions occur. These become important for energies above the Coulomb barrier  $E_c$  (in MeV):

$$E_c \approx Z_1 Z_2 (A_1^{1/3} + A_2^{1/3}) \quad (4)$$

When during nuclear reactions the irradiative particle is absorbed by the nucleus, the target atom experiences a recoil energy which is  $E_r \cdot f$ . When the capture is followed by emission of another particle with mass  $M_3 \approx M_1 \approx M_2$  and energy  $E_3$  (in the center of mass system), then the recoil energy of the daughter nucleus is in the interval:

$$T_r = (T_{max}/4) (1 + E_3/E) \approx 2\sqrt{E_3/E} \quad (5)$$



1. Read fundamentals
2. Search for data about high-energy precision irradiation
3. Check the response of biological material in relation to the applied dose



**Thank you !**  
**Stay tuned !**

# Backup Slides

<i>user:</i>	researcher, engineer
<i>field of activity:</i>	medical research and application
<i>sample task:</i>	design of appropriate devices and beam lines within an tumor treatment installation
<i>SpringerMaterials usage:</i>	type in easy search box: "ion irradiation"
<i>result:</i>	first hit: overview of the physical concepts and relevant quantities with respect to prominent particles.
<i>volume:</i>	LB III/25

The information about ion irradiation enables any engineer or medical physicist to design and implement appropriate accelerators and beam lines. For that he has to know the energy loss and damage rates of protons, for example, as described in vol. LB III/25. With 'ion irradiation' typed into the easy-search field he obtains the needed information.

<i>user:</i>	scientist
<i>field of activity:</i>	chemical analysis
<i>sample task:</i>	measuring by NMR spectroscopy the chemical shift of the material to be analyzed and comparing it with the chemical shift of the expected material $C_9H_{14}N_2$
<i>SpringerMaterials usage:</i>	type in "C9H14N2 chemical shift" into easy-search box; select "Chemical shifts and coupling constants of C9H14N2"
<i>result:</i>	hit: " $C_9H_{14}N_2$ " contains the desired data
<i>volume:</i>	LB III/35D3

A chemical analyst shall find out if the material to be analyzed is the expected material  $C_9H_{14}N_2$ . He uses NMR spectroscopy for this task and likes to compare the measured chemical shift with the chemical shift of  $C_9H_{14}N_2$ . For this he types "C9H14N2 chemical shift" into the easy-search box and selects "Chemical shifts and coupling constants of  $C_9H_{14}N_2$ ", the hit "C9H14N2" contains the desired data.

<i>user:</i>	researcher, engineer
<i>field of activity:</i>	sensor development
<i>general task:</i>	design of a flame sensor
<i>sample task 1:</i>	get an overview on pyroelectricity
<i>SpringerMaterials usage:</i>	type "pyroelectric effect" into easy-search box
<i>result:</i>	3rd hit: "Pyroelectric equations of state and definition of pyroelectric coefficients"
<i>volume:</i>	LB III/29B
<i>sample task 2:</i>	get suitable materials
<i>SpringerMaterials usage:</i>	type "pyroelectric coefficient" into Advanced Search Properties box

~~A researcher or an engineer in a department for sensor development wants to design a new flame sensor. First, he might wish to refresh his knowledge on pyroelectricity. He types "pyroelectric effect" in the easy search. The third hit leads him to a chapter on fundamentals in volume III/29B. After reading, he wants to find suitable materials, for which he goes to the Advanced Search and types "pyroelectric coefficient" into the Properties search field. He receives a lot of documents about relevant materials and will then be able to choose the best by comparing their pyroelectric coefficients.~~

91 hits in electronic "Structure Structure and Transport", mostly substances  
LB III/29B, III/36B

<i>user:</i>	researcher, engineer
<i>field of activity:</i>	permanent magnet development
<i>sample task:</i>	get an overview of materials with magnetic hysteresis
<i>SpringerMaterials</i>	type "magnetic hysteresis" into easy-search box
<i>usage:</i>	
<i>result:</i>	51 hits in "Magnetism"
<i>volume:</i>	LB III/27, III/19

A company wants to develop permanent magnets for special applications defined by the company's customers. A researcher/engineer of the development department gets the job to gather information for suitable materials. He/she types "magnetic hysteresis" into the easy-search box and gets a list of materials which show magnetic hysteresis. This list may be used to continue search for special materials suitable for the customer's needs.

<i>user:</i>	researcher, engineer
<i>field of activity:</i>	Photovoltaics
<i>task:</i>	<ul style="list-style-type: none"> <li>• design of a solar cell prototype</li> <li>• gain of overview on the state of knowledge of the underlying physical concepts</li> <li>• test of a suitable new material; information about this material's properties</li> </ul>
<i>SpringerMaterials usage:</i>	type in "photovoltaic cell" into easy-search box; select "photovoltaic power generation"
<i>result:</i>	first hit: overview of current established state of concepts and techniques around solar cells ; selection of $\text{CuGaSe}_2$ as material of interest
<i>SpringerMaterials usage:</i>	type in "CuGaSe2" into substance-box under 'Advanced Search', "band energy" into property-box
<i>result:</i>	first hit: data around $\text{CuGaSe}_2$ needed to project a solar device (e.g. suitability to absorb light dependent on temperature, cf. Fig. 3)
<i>volume:</i>	LB III/41E
<i>SpringerMaterials usage:</i>	alternatively type in "CuGaSe2" into substance-box under 'Advanced Search', "mobility" into property-box
<i>result:</i>	first hit: data of the charge carriers of $\text{CuGaSe}_2$ , in particular mobility and concentration
<i>volume:</i>	LB III/41E

A solar technician, employed to find a material system for a new photovoltaic cell, might type 'photovoltaic cell' simply into the SpringerMaterials easy-search template, resulting in a first hit in a document about 'photovoltaic power generation'. From that he gains a comprehensive overview about the concepts and techniques around solar cells. He might learn from the document that  $\text{CuGaSe}_2$  may be a suitable material of choice for his intentions. The amount of the energy gap and particularly its behavior must therefore be known. He uses the 'Advanced Search', typing ' $\text{CuGaSe}_2$ ' into the substance-field along with 'band energy' into the property-field. As a result SpringerMaterials offers as first hit a document that describes the data needed in the context of different other properties. Fig. 3 even displays the dependence of the band gap energy from the temperature.

He might be even interested in the carrier mobility within  $\text{CuGaSe}_2$ , typing in 'mobility' under 'property', thus ending up at a document about transport properties of that material with all relevant quantities.



<i>user:</i>	researcher, engineer
<i>field of activity:</i>	accelerator design
<i>task:</i>	develop a superconducting magnet
<i>task:</i>	get the properties of wires of Nb <sub>3</sub> Sn
<i>SpringerMaterials usage:</i>	type "Nb3Sn wire" into easy-search box
<i>result:</i>	26 hits in "Superconductivity"
<i>volume:</i>	LB III/21

A research institute is designing a new particle accelerator which allows to accelerate particles like protons or heavy ions to increasingly higher energy before they hit each other. This can be achieved in circular accelerators provided the magnets used to bent the particle beam sustain a sufficiently high magnetic field. It is known that superconducting Nb<sub>3</sub>Sn would be a suitable material. As the magnet would be made of wires the researcher/engineer types "Nb3Sn wire" into the easy-search box to get documents which provide information on the properties of wires of Nb<sub>3</sub>Sn.

<i>user:</i>	engineer
<i>field of activity:</i>	electronic engineering industry
<i>task:</i>	miniaturization of storage elements
<i>SpringerMaterials usage:</i>	type in easy-search box: "Ba-ferrite magnetic recording"
<i>result:</i>	first hit: overview and suitability of a series of Ba-ferrites in the realm of magnetic recording is presented, along with properties
<i>volume:</i>	LB III/27G
<i>SpringerMaterials usage:</i>	'Advanced Search', type in substance-box: "Ba-ferrite", in property-box "crystal structure"
<i>result:</i>	first hits: data on the crystal structure of Ba-ferrites with further references
<i>volume:</i>	LB III/7 (possibly f)
<i>result:</i>	when selecting 'Inorganic Solid Phases' (LPDF-database) even more precise data are accessible

*volume:* LPF Database

An electronic engineer might strive to enhance the storage capacity of hard disks. To minimize its spatial dimensions the magnetic behavior of hexagonal ferrites are of interest, for instance barium-ferrite. When typing 'Ba-ferrite' into the 'substance'-field and 'crystal structure' into the property-field he obtains as first hit information about the crystal structure and its behavior under different circumstances. When selecting 'Inorganic Solid Phases' he is offered even more precise data. When typing 'Ba-ferrite' along with 'magnetic recording' an overview of hexagonal ferrites is displayed as first hit, that are of interest as possible recording devices.

<i>user:</i>	researcher, scientist
<i>field of activity:</i>	scientific investigation of impurities within a semiconductor
<i>task:</i>	design of an annealing-installation for semiconductor after implantation of atoms under inert gas flow; choice: Argon gas mixture
<i>SpringerMaterials usage:</i>	under 'Advanced Search' type in "Argon" into the substance-box and "virial coefficient" into the property-box
<i>result:</i>	first hit: overview of several gas mixtures containing Ar with the corresponding virial coefficients
<i>volume:</i>	LB IV/21B

For any engineer or scientist in a laboratory concerned with the production of technical material, a major means to alter the material's microstructure is annealing. In order to avoid unintentional absorption of foreign material from the surrounding ('extrinsic defects') the process takes place in inert gas atmosphere. Argon, and mixtures with other gases, is a widely used representative gas.

The knowledge of the virial coefficients allows the exact calculation of the gas pressure, for example, if a certain temperature is required and volume given.

By using the 'Advanced Search' and typing 'Argon' into the substance-field, 'virial coefficient' into the property-field, the user gets as first hit a list of these properties for several mixtures of Argon (LB IV/21B).

<i>user:</i>	researcher, engineer
<i>field of activity:</i>	health risk assessment
<i>task:</i>	get information on Pu alloys and compounds
<i>task:</i>	get all available phase diagrams of compounds containing Pu
<i>SpringerMaterials usage:</i>	type "plutonium phase diagram" into easy-search box
<i>result:</i>	113 hits in "Multiphase Systems"
<i>volume:</i>	LB IV/5, IV/11

For a nuclear power plant to be built a national health institute or engineering company has to assess the health risk to the employees of the power plant and the people living in its neighborhood that would result from a malfunctioning of the power plant. To this end knowledge of the properties of plutonium (one of the most hazardous materials) and plutonium-containing compounds is of paramount importance. Phase diagrams of plutonium compounds allow to determine which compounds might possibly be generated under the conditions of operation of the power plant. Therefore the researcher/engineer types "plutonium phase diagram" into the easy-search box. Similarly the phase diagrams of other hazardous materials may be found.

<i>user:</i>	researcher, engineer
<i>field of activity:</i>	displays
<i>task:</i>	reducing the switching time of a liquid crystal display
<i>SpringerMaterials usage:</i>	type in "liquid crystal switching time" into easy-search box
<i>result:</i>	first hit: "liquid crystals - definition of properties" – An overview on all properties is given here. Search "switching time" in this document. – The researcher finds out: "It is desirable to have substances with low viscosity for application in a LC cell."
<i>volume:</i>	LB VIII/5A
<i>task:</i>	Now the researcher likes to get viscosity data of liquid crystals.
<i>SpringerMaterials usage:</i>	type in "liquid crystal viscosity" into easy-search box.
<i>result:</i>	first hit: "liquid crystals - definition of properties" (as above), second hit: "liquid crystals – introduction" – In this file the arrangement of the data files is explained. hit 8: "Data ordered by properties > Viscosity" contains the desired data

A researcher has the task to reduce the switching time of a liquid crystal display. He first likes to find out on which properties the switching time depends and types in "liquid crystal switching time" into the easy-search box. The first hit is a document giving an overview over all properties of liquid crystals. By searching "switching time" in this document the researcher finds out that it is desirable to have substances with low viscosity. He now searches for "liquid crystal viscosity" and finds an introduction where the arrangement of the data files on liquid crystals is explained and a data file "Data ordered by properties – Viscosity" which contains the desired viscosity data.

<i>user:</i>	scientist, engineer
<i>field of activity:</i>	laser processing
<i>task:</i>	- metal sheet cutting - gain of overview on the state of knowledge
<i>SpringerMaterials usage:</i>	type in "cutting" into easy-search box; select "cutting"
<i>result:</i>	first hit: "Cutting: Modeling and data" - overview on the state of knowledge
<i>volume:</i>	LB VIII/1C

A scientist/an engineer working in the field of laser processing likes to gain an overview on metal sheet cutting. He simply types in “cutting” into the easy-search box. The first hit is the desired overview article “Cutting: Modeling and data”.

<i>user:</i>	astronomer, researcher, engineer
<i>field of activity:</i>	observatory
<i>task:</i>	find information about observed asteroid
<i>SpringerMaterials usage:</i>	type in "asteroid" into easy-search box; select "asteroid"
<i>result:</i>	first hit: "Astronomy and Astrophysics before 1993" is not the correct one, as mainly actual information is searched for; second hit: "The asteroids" gives an overview on asteroids
<i>volume:</i>	LB VI/4B

An astronomer or a scientist interested in astronomy who has observed an asteroid wants to find information about asteroids. In Landolt-Börnstein an overview on asteroids is given.

<i>user:</i>	scientist, researcher, engineer
<i>field of activity:</i>	researcher in optics industry
<i>task:</i>	design of an infrared detector
<i>SpringerMaterials usage:</i>	type in 'Advanced Search' into substance-box "InSb" and "energy gap" into property-box; select LPF Database by marking the corresponding box to the left
<i>result:</i>	first hit yields all LPF-data sheets on InSb containing physical property; click first hit; seventh hit from top: overview of the physical concepts and relevant quantities with respect to prominent particles.
<i>volume:</i>	LPF Database

A researcher in the optics industry wants to design a new infrared detector. He knows that an indium antimony alloy is a possible material and needs to find data about the energy gap of this material. E.g., he uses the substance-property search in the Advanced Search section. The first hit leads to the summary document of the InSb entry in the LPF database, then the seventh hit from top contains the data on the energy gap.



<i>user:</i>	researcher, engineer
<i>field of activity:</i>	chemical engineering
<i>task:</i>	calculate the freezing point depression of ethanol/water mixtures
<i>task:</i>	get the thermophysical properties of ethanol/water mixtures
<i>SpringerMaterials</i>	choose "Thermophysical Properties", type "ethanol" and
<i>usage:</i>	choose "ethanol + water"
<i>result:</i>	1 hit in "Thermophysical Properties"
<i>volume:</i>	Dortmund Data Bank

A car manufacturer wants to provide information in the cars' manuals on how much ethanol is to be used in ethanol/water mixtures so that the mixture does not freeze even in strong winters and can be still applied for cleaning the window panes.

To give such an advice the freezing point of ethanol/water mixtures has to be calculated in dependence on the ethanol content. The researcher/engineer types "ethanol" and chooses "ethanol + water" in the input box of the Thermophysical Properties subject area.

From the thermophysical data provided the depression of the freezing point by adding ethanol to water can be calculated by the researcher/engineer.

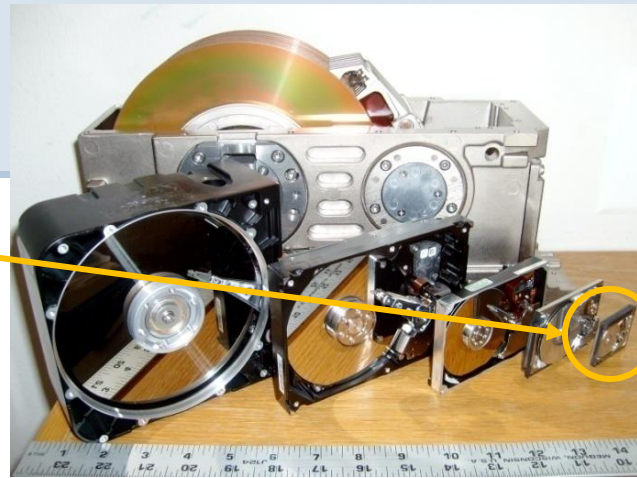
## Inorganic Solid Phases

- The database on Inorganic Solid Phases contains all necessary data for any researcher or scientist dealing with materials in their solid form. It provides information about the regions and conditions of their existence (phase diagrams), their crystallographic and geometric structure (structural data) and their physical properties (property data).

## Thermophysical Properties

- Thermophysical data of pure organic substances as well as mixtures consisting of an organic substance and water, or two organic substances, are represented in tables and diagrams. Chemical engineers are able to design and develop chemical reactors and plants if they know the thermophysical properties of the substances used in these reactors and plants. The data is also useful for (organic) chemists in research laboratories.

Engineer develops smaller hard disk with higher memory capacity



# SpringerMaterials

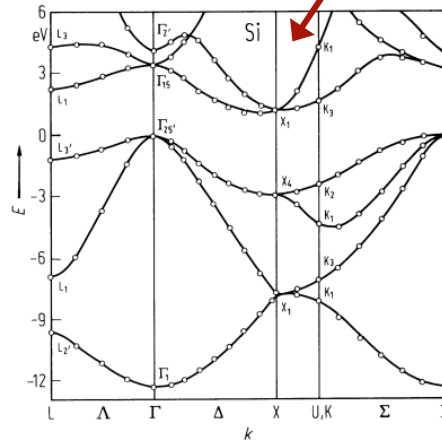
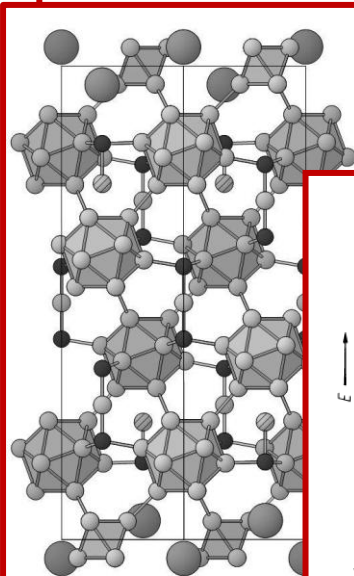
## Typical Usage Situation

Ref. p. 216] 6.3.11 Magnetic recording applications of hexagonal ferrites 209

### 6.3.11 Magnetic recording applications of hexagonal ferrites

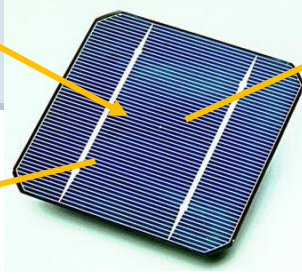
Table 32.

Device	Ref.	Remarks	Fig./Table
recording (general)	85F1	survey of applications of Ba-ferrite media to perpendicular recording devices	
recording (general)	85F2	review on the potential application of Ba-ferrite media for perpendicular magnetic recording	123, 124
recording (general)	85H	development of Ba-ferrite particulate media for perpendicular recording applications	
recording (general)	85M1	Ba-ferrite media for perpendicular recording, recording density vs. degree of orientation, thermal stability	
recording (general)	85V3	development of Ba-ferrite particulate media for application to VCRs, VTRs, DATs and high density floppy disks	
recording (general)	86I1	achievements in magnetic recording applications of Ba-ferrite in Japan	
recording (general)	87F1	Ba-ferrite particulate media for perpendicular recording, output level vs. particle aspect ratio and size distribution	125
recording (general)	87P	dynamic model for simulating the magnetic recording processes, application to Ba-ferrite media	
recording (general)	87S2	recording characteristics of oriented and non-oriented Ba-ferrite media	
recording (general)	87S3	comparative analysis of particulate media for magnetic recording vs. particle and coating characteristics	
recording (general)	88S7	comparison between Ba-ferrite and Co- $\gamma$ -Fe <sub>2</sub> O <sub>3</sub> films (0.25 $\mu$ m thick), Ba-ferrite media show the highest linear density ( $> 1400$ kbit/mm) and a very sharp distribution of remanent coercivity	
recording (general)	88S3	analysis of possible solutions to a rewrite modulation in Ba-ferrite particulate media	
recording (general)	89V2	application of dynamic micromagnetic theory to signal and noise problem in Ba-ferrite media	
developer for electro-photography duplicator for R-DAT	83	Ba(Ni,Zn)-Y ferrite as carrier for dual component developer	
	86N	Ba-ferrite media for magnetic contact duplication at $\lambda$ from 0.6 to 40 $\mu$ m, track width from 10 to 20 $\mu$ m, block error rate = $5 \cdot 10^{-7}$	
Floppy disk	85P3	Ba-ferrite perpendicular recording media, 3.5" FD, endurance test	
Floppy disk	85I1	Ba-ferrite perpendicular recording media, 3.5" 4 MB FD, recording density = 35 kbit, track density = 135 TPI	Table K
Floppy disk	85I2	Ba-ferrite perpendicular recording media, flexible disks for electronic still cameras ( $\phi = 47$ mm, 7400 rpm)	128
Floppy disk	86V	Ba-ferrite media (coating thickness 3 $\mu$ m, $\rho_2 = 0.54$ ), high density recording response and overwrite characteristics	



1. Search suited magnetic material
2. Get information about its molecular structure
3. Check its physical properties

Engineer develops improved solar cell



# SpringerMaterials

## Typical Usage Situation

1. Search for the carrier mobility of CuGaSe<sub>2</sub>
2. Get an overview of the climate conditions in the area of interest

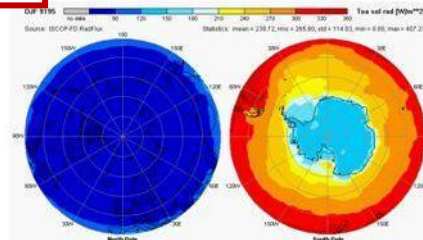
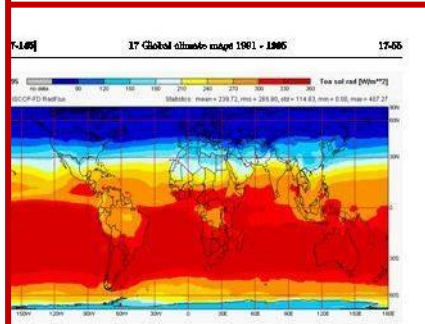
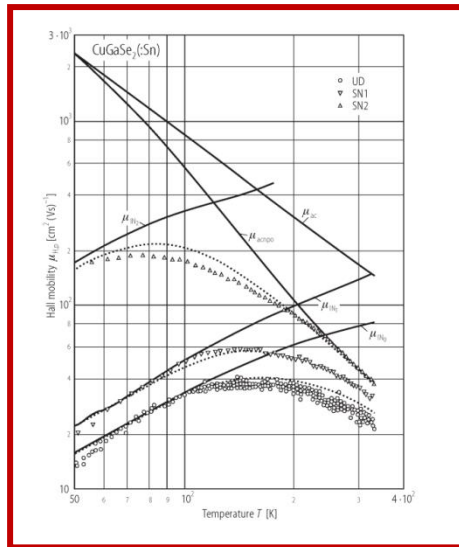


Fig. 17.46. IPCC - mean winter and solar radiation flux density across top of atmosphere (1961-1996). Solar radiation is positive throughout (downwards).

Engineer needs to improve the corrosion resistance of car bodies:  
Steels coated with Al-Fe-Zn layer



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## Typical Usage Situation

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Go Advanced Search

Home Bookshelf Periodic Table Help For Librarians Feedback

Al-Fe-Zn

Al-Fe-O-Zn

Al-Ba-Fe-O-Zn

Al-Co-Fe-O-Zn

Al-Cr-Fe-O-Zn

Al-Fe-Li-O-Zn

Al-Fe-Mg-O-Zn

Al-Fe-Mn-Si-Zn

Al-Ba-Ca-Fe-O-Zn

Al-Ba-Fe-La-O-Zn

Al-Be-Fe-Mg-O-Zn

Al-Co-Cr-Fe-O-Zn

Al-Cr-Fe-Mg-O-Zn

Al-Cr-Fe-Ni-O-Zn

Al-Fe-H-O-S-Zn

Al-Fe-H-O-Sn-Zn

Al-Fe-Mg-Mn-O-Zn

Al-Fe-Nd-O-Ti-Zn

Al-Cr-Fe-Mg-O-Ti-Zn

Al-Fe-H-K-O-S-Zn

Al-Fe-H-Mg-O-Sn-Zn

Al-Fe-H-Mg-O-Ti-Zn

Al-Fe-H-O-S-Ti-Zn

Al-Fe-Li-Mn-O-Ti-Zn

Al-Fe-Mn-O-Pb-Ti-Zn

Al-Be-F-Fe-H-Mg-O-Zn

Al-Ca-Fe-Mg-Mn-O-Si-Zn

Al-Cr-Fe-H-Mg-O-Ti-Zn

Search for Element Systems

Select elements by clicking on the symbols.  
Deselect elements by clicking a second time.

Al-Fe-Zn

1. Select components by clicking on elements in periodic table
2. Available content shown

The Landolt-Börnstein Database

Go Advanced Search

Periodic Table Help For Librarians Feedback

Results 1 - 7 of 7 Documents

Compact View Clear Refine

Al-Fe-Zn

Al-Fe-Zn, ternary phase diagram, isothermal section

Al-Fe-Zn, ternary phase diagram, liquidus projection

Al-Fe-Zn, ternary phase diagram, vertical section

Tracer diffusion coefficients for ternary alloys

Al-Fe-Zn, ternary phase diagram, isothermal section

Al-Fe-Zn, ternary phase diagram, liquidus projection

Al-Fe-Zn, ternary phase diagram, vertical section

Tracer diffusion coefficients for ternary alloys

Al-Fe-Zn, ternary phase diagram, isothermal section

Al-Fe-Zn, ternary phase diagram, liquidus projection

Al-Fe-Zn, ternary phase diagram, vertical section

Tracer diffusion coefficients for ternary alloys

Engineer needs to improve the corrosion resistance of car bodies: Steels coated with Al-Fe-Zn layer



# SpringerMaterials Typical Usage Situation

Substances, Properties, ... Bibliographic References Help Close

Your Query  
"Al-Fe-Zn" automotive Go Clear

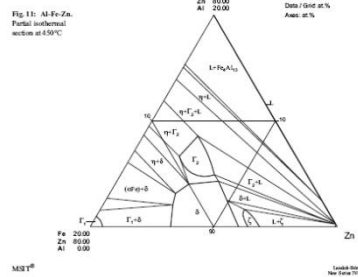
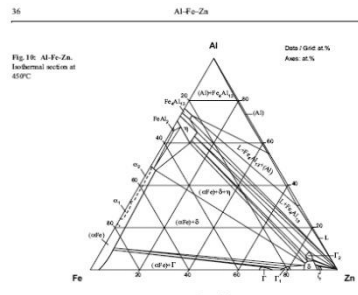
Search for ...  
Substances / Molecular Formulas / Element Systems / CAS Registry Numbers

Properties

Search in ... Search for ...

- Particles, Nuclei and Atoms
- Molecules and Radicals
- Electronic Structure and Transport
- Magnetism
- Semiconductivity
- Superconductivity
- Crystallography
- Thermodynamics
- Multiphase Systems
- Advanced Materials
- Advanced Technologies
- Astro- and Geophysics

Inorganic Solid Phases  
 Thermophysical Properties  
 Chemical Safety



Phase / Temperature Range [°C]	Pearson Symbol / Space Group / Prototype	Lattice Parameters [nm]	Comments/References
FeAl <sub>2</sub> < 1156	aP18 P1 FeAl <sub>2</sub>	a = 487.3 b = 466.1 c = 860.0 α = 91.75° β = 73.27° γ = 96.89°	[2001Pa], at 66.9 at.% Al solid solubility ranges from 65.5 to 67.0 at.% Al
η-FeAl <sub>3</sub> < 1169	cC24 Cmcm	a = 765.59 b = 641.54 c = 421.84	[2001Pa], at 71.5 at.% Al solid solubility ranges from 71.0 to 72.5 at.% Al Equilibrium solubility is up to 11 at.% Zn at 450°C [1992Pe] [2001Ko], at FeAl <sub>3</sub> Zn
FeAl <sub>11</sub> < 1160	hC102 C2/m FeAl <sub>11</sub>	a = 1552.7 to 1568.7 b = 803.5 to 808.4 c = 1244.9 to 1248.8 β = 107.7 to 107.99°	[2001Pa], 74.16 to 76.7 at.% Al solid solubility ranges from 74.5 to 75.5 at.% Al [2001Pa], at 76.0 at.% Al
FeAl <sub>12</sub> < 1160	hC102 C2/m FeAl <sub>12</sub>	a = 1549.2 b = 807.8 c = 1247.1 β = 107.60°	sometimes called FeAl <sub>3</sub> in the literature
FeAl <sub>13</sub> < 1160	cP28 Fm-3m FeAl <sub>13</sub>	a = 897.41 b = 901.8	[V-C], solid solubility ranges from 68.0 to 82.5 at.% Zn
FeAl <sub>14</sub> < 1160	cF208 Fm-3m FeAl <sub>14</sub>	a = 1796.3	[V-C2], solid solubility ranges from 75.5 to 81.0 at.% Zn
FeAl <sub>15</sub> < 1160	hC102 C2/m FeAl <sub>15</sub>	a = 1283.0 b = 3770.0	[V-C], solid solubility ranges from 86.5 to 92.0 at.% Zn Equilibrium solubility is up to 4.3 at.% Al at 450°C [1992Pe]
FeAl <sub>16</sub> < 1160	cF208 Fm-3m FeAl <sub>16</sub>	a = 1342.4 b = 760.8 c = 506.1 β = 127.3°	[V-C], solid solubility ranges from 92.5 to 96.0 at.% Zn Equilibrium solubility is up to 1.05 at.% Al at 450°C [1992Pe]
FeAl <sub>17</sub> < 1160	-	-	[1992Pe, 1989Van]

- Refine search and check available content
- Check phase diagram and crystallographic data of the chosen composition

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"Al-Fe-Zn" automotive Go Advanced Search

Home Bookshelf Periodic Table Help For Librarians Feedback

0 Particles, Nuclei and Atoms  
0 Molecules and Radicals  
0 Electronic Structure and Transport  
0 Magnetism  
0 Semiconductivity  
0 Superconductivity  
0 Crystallography  
0 Thermodynamics  
2 Multiphase Systems  
0 Advanced Materials  
0 Advanced Technologies  
0 Astro- and Geophysics  
0 Inorganic Solid Phases  
0 Thermophysical Properties  
0 Chemical Safety

Results 1 - 2 of 2 Documents Compact view Clear Refine

Multiphase Systems > Ternary Alloys > Phase Diagrams, Crystallography and Thermodynamics > Light Metal Systems > Aluminum (Al-X-Y) Ternary Alloys  
Al-Fe-Zn

Metadata - Substance: Al-Fe-Zn ... Metadata - Element System: Al-Fe-Zn ... Fulltext: steel sheets for the automotive industry. As a result, recent ...

Multiphase Systems > Ternary Alloys > Phase Diagrams, Crystallography and Thermodynamics > Iron Systems > Selected Systems from Al-B-Fe to C-Cu-Fe  
Aluminum - Iron - Zinc

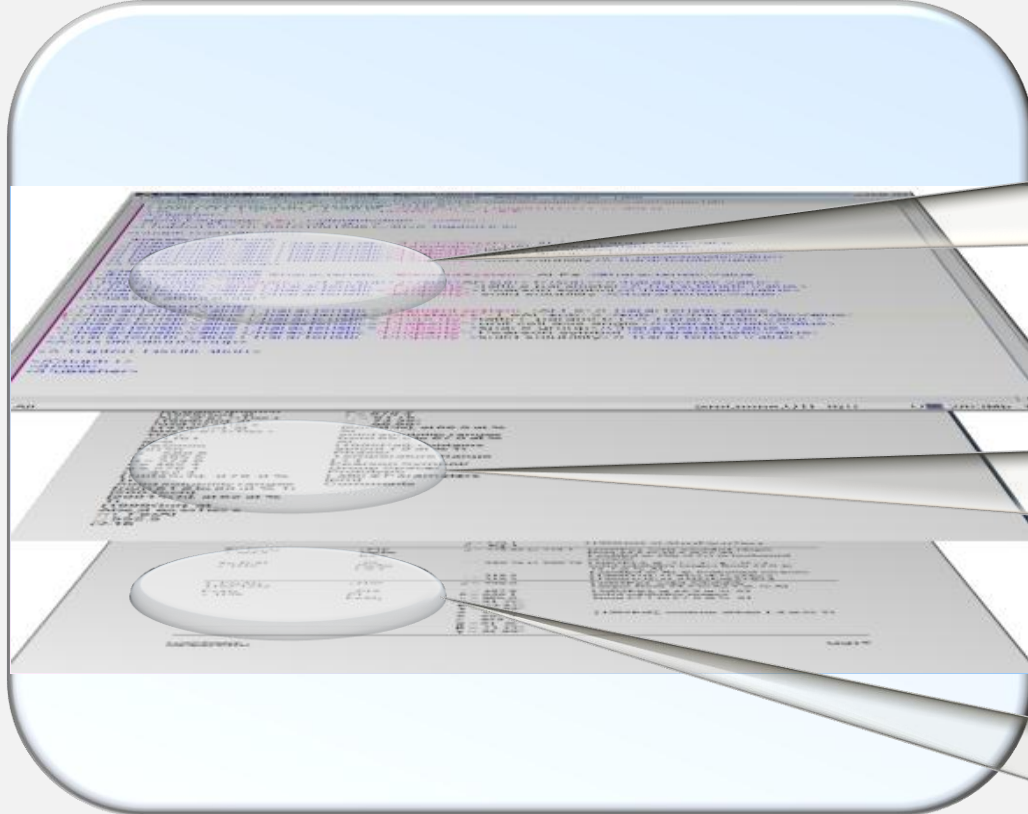
Metadata - Substance: Al-Fe-Zn ... Metadata - Element System: Al-Fe-Zn ... Fulltext: steel sheets for the automotive industry. As a result, recent ...

Three Layers per Doc

A Closer Look

Usability

# Summary



```

<ClassificationGroup>
<CharacteristicValue Characteristic='ElementSystem'>Al-Fe</CharacteristicValue>
<CharacteristicValue Characteristic='Substance'>FeAl<sub>2</sub></CharacteristicValue>
<CharacteristicValue Characteristic='Property'>unit cell axis angle</CharacteristicValue>
<CharacteristicValue Characteristic='Property'>Pearson symbol</CharacteristicValue>
<CharacteristicValue Characteristic='Property'>solid solubility</CharacteristicValue>
</ClassificationGroup>
  
```

**precise search result**

Ti.  $aP18$   
 Solid solubility ranges from 33.5 to 53.3 at.%  
 Ti [2003Sch].  $P1$   
 [2003Sch], at 38 at.%  $FeAl_2$   
 Ti.  $a = 487.8$   
 [2000Mab], at Al-47  $b = 646.1$   
 at.% Ti.  $c = 880.0$   
 Heat treated at 1000°C  $\alpha = 91.75^\circ$   
 for 48 h followed by water quench.  $\beta = 73.27^\circ$   
 [1999Gor] at  $\gamma = 96.89^\circ$   
 $c = 879.4$

**wrong search result**

$FeAl_2$	$aP18$	$a = 487.8$	[2003Pis], at 66
$\leq 1156$	$P1$	$b = 646.1$	solid solubility
	$FeAl_2$	$c = 880.0$	from 65.5 to 67.
		$\alpha = 91.75^\circ$	
		$\beta = 73.27^\circ$	
		$\gamma = 96.89^\circ$	
		$c = 879.4$	
		$\alpha = 91.76^\circ$	[1995Pal], cont
		$\beta = 73.35^\circ$	
		$\gamma = 96.89^\circ$	

**easy to understand but not searchable**

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What the user can find with the SpringerMaterials Search engine

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What the user can find with the SpringerLink search engine



What the user sees

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