

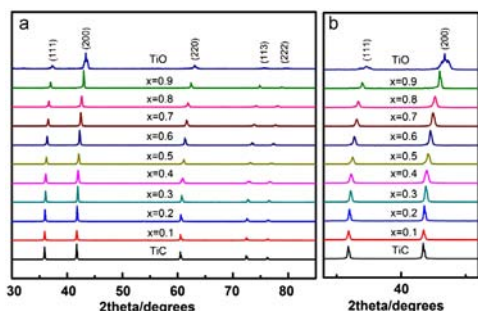
Abstracted/indexed in BioEngineering Abstracts, Chemical Abstracts, Coal Abstracts, Current Contents/Physics, Chemical, & Earth Sciences, Engineering Index, Research Alert, SCISEARCH, Science Abstracts, and Science Citation Index. Also covered in the abstract and citation database SciVerse SCOPUS<sup>®</sup>. Full text available on SciVerse ScienceDirect<sup>®</sup>.

## Regular Articles

### Structural studies of $\text{TiC}_{1-x}\text{O}_x$ solid solution by Rietveld refinement and first-principles calculations

Bo Jiang, Na Hou, Shanyan Huang, Gege Zhou, Jungang Hou, Zhanmin Cao and Hongmin Zhu

page 1



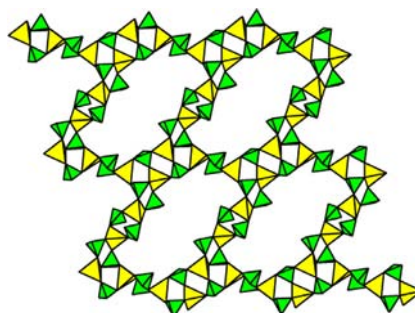
XRD of series of titanium oxycarbides ( $\text{TiC}_{1-x}\text{O}_x$ ,  $0 \leq x \leq 1$ ) solid solution prepared by adjusting the proportion of TiO in the starting material.

## Regular Articles—Continued

### Synthesis and characterization of three-layered zinc phosphites containing *tert*-octylamine molecules with template and ligand roles

Chih-Min Wang, Cheng-Wei Chiu, Hsiu-Mei Lin and Kwang-Hwa Lii

page 16

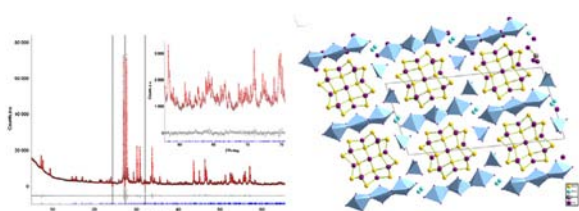


The first layered metal phosphite with large 20-ring windows,  $(\text{C}_8\text{H}_{17}\text{NH}_3)_4\text{Zn}_3(\text{HPO}_3)_5 \cdot 3\text{H}_2\text{O}$  (1), sandwiched by water molecules and protonated *tert*-octylamine molecules.

### Cobalt-doped $\text{Bi}_{26}\text{Mo}_{10}\text{O}_{69}$ : Crystal structure and conductivity

Z.A. Mikhailovskaya, E.S. Buyanova, S.A. Petrova, M.V. Morozova, V.M. Zhukovskiy, R.G. Zakharov, N.V. Tarakina and I.F. Berger

page 9

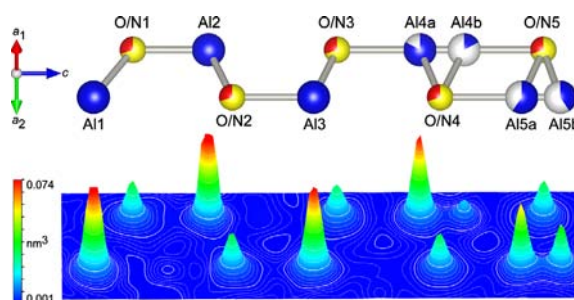


Measured and calculated diffraction spectra for  $\text{Bi}_{12.8}\text{Co}_{0.2}\text{Mo}_5\text{O}_{34 \pm \delta}$  and projection of the  $\text{Bi}_{12.8}\text{Co}_{0.2}\text{Mo}_5\text{O}_{34 \pm \delta}$  crystal structure onto the *ac* plane.

### Electron density distribution and crystal structure of 27*R*- $\text{Al}_9\text{O}_3\text{N}_7$

Toru Asaka, Hiroki Banno, Shiro Funahashi, Naoto Hirosaki and Koichiro Fukuda

page 21



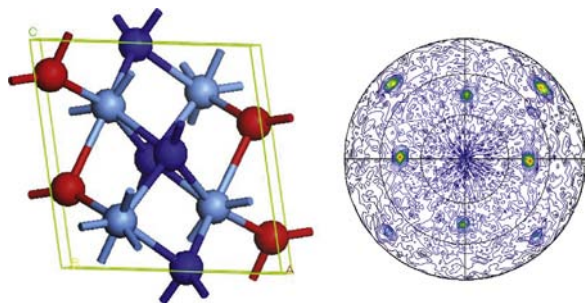
A bird's eye view of electron densities up to 50% ( $0.074 \text{ nm}^{-3}$ ) of the maximum on the plane parallel to (110) with the corresponding atomic arrangements of  $\text{Al}_9\text{O}_3\text{N}_7$ .

Continued

### Growth of single crystalline TaON on yttria-stabilized zirconia (YSZ)

Junguang Tao, J.W. Chai, L.M. Wong, Z. Zhang, J.S. Pan and S.J. Wang

page 27

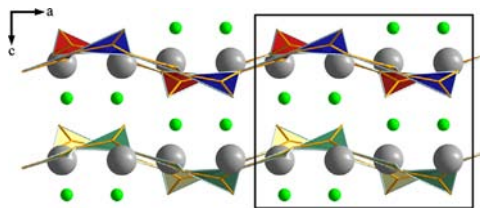


Structure of single crystalline  $\beta$ -TaON and its diffraction pole figure.

### High-pressure synthesis and characterization of the first cerium fluoride borate $\text{CeB}_2\text{O}_4\text{F}$

Ernst Hinteregger, Klaus Wurst, Martina Tribus and Hubert Huppertz

page 47

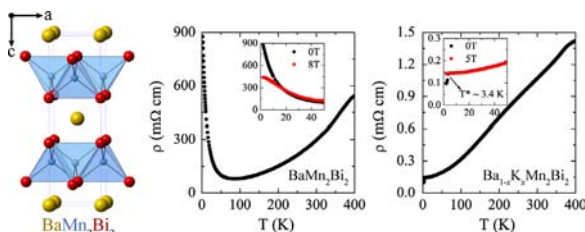


A new rare-earth fluoride borate  $\text{CeB}_2\text{O}_4\text{F}$  could be synthesized under high-pressure/high-temperature conditions of 0.9 °GPa and 1450 °C in a Walker-type multianvil apparatus. The crystal structure represents a new structure type in the class of rare-earth fluoride borates. The structure exhibits a 9+1 coordinated cerium ion, one three-fold coordinated fluoride ion and a one-dimensional chain of  $[\text{BO}_3]^{3-}$  groups. A closer view on the  $ac$ -plane shows an interesting wave-like modulation of the borate chains.

### Crystals, magnetic and electronic properties of a new $\text{ThCr}_2\text{Si}_2$ -type $\text{BaMn}_2\text{Bi}_2$ and K-doped compositions

Bayrammurad Saparov and Athena S. Sefat

page 32

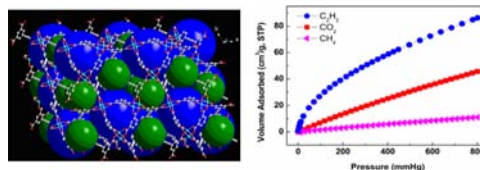


Local moment antiferromagnet  $\text{BaMn}_2\text{Bi}_2$ , the first bismuthide with  $\text{ThCr}_2\text{Si}_2$  structure, turns metallic upon K-doping.

### A microporous metal–organic framework with butynylene functionality for selective gas sorption

Lifeng Wang, Lu Zhai, Xiaoming Ren and Wenwei Zhang

page 53

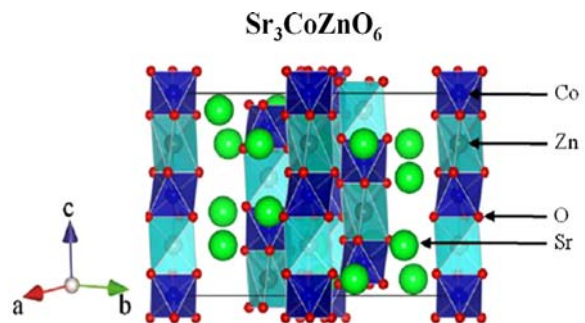


A microporous metal–organic framework with butynylene groups,  $\{[\text{Cu}_2(\text{BBTC})(\text{H}_2\text{O})_2] \cdot 2\text{DMSO} \cdot 4\text{H}_2\text{O}\}_n$ , exhibits higher gas selective sorption towards  $\text{C}_2\text{H}_2/\text{CH}_4$  and  $\text{CO}_2/\text{CH}_4$  at room temperature with a selectivity of 5.7 and 4.1, respectively.

### Crystal structure and magnetic properties and Zn substitution effects on the spin-chain compound $\text{Sr}_3\text{Co}_2\text{O}_6$

Xia Wang, Yanfeng Guo, Ying Sun, Yoshihiro Tsujimoto, Yoshitaka Matsushita and Kazunari Yamaura

page 40

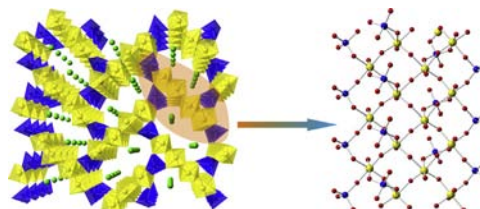


Crystal structure of the spin-chain compound  $\text{Sr}_3\text{CoZnO}_6$  synthesized at 6 GPa. Zn atoms preferably occupy the trigonal prism sites rather than the octahedral sites. As a result, the compound is much magnetically isotropic.

### $\text{K}[\text{AsW}_2\text{O}_9]$ , the first member of the arsenate–tungsten bronze family: Synthesis, structure, spectroscopic and non-linear optical properties

Evgeny V. Alekseev, Olivier Felbinger, Shijun Wu, Thomas Malcherek, Wulf Depmeier, Giuseppe Modolo, Thorsten M. Gesing, Sergey V. Krivovichev, Evgeny V. Suleimanov, Tatiana A. Gavrilova, Lev D. Pokrovsky, Alexey M. Pugachev, Nikolay V. Surovtsev and Victor V. Atuchin

page 59



$\text{K}[\text{AsW}_2\text{O}_9]$ , the first member of arsenate–tungsten bronze family exhibit new three dimensional structure type, significant thermal stability and NLO properties.

Download English Version:

<https://daneshyari.com/en/article/7759787>

Download Persian Version:

<https://daneshyari.com/article/7759787>

[Daneshyari.com](https://daneshyari.com)