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CONTENTS

Foreword

Professor Hans Georg von Schnering celebrates his 75th birthday

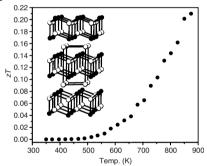
Reinhard Nesper and Yuri Grin *Page 2251*

Special Section:

Dedicated to the occasion of the 75th birthday of Prof. Hans Georg von Schnering

Thermoelectric properties and microstructure of Mg₃Sb₂ Cathie L. Condron, Susan M. Kauzlarich, Franck Gascoin and G. Jeffrey Snyder

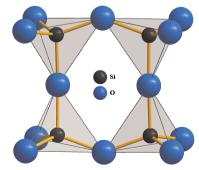
Page 2252



Dimensionless figure of merit for Mg_3Sb_2 hot pressed and sintered at 873 K. The inset illustrates the crystal structure of Mg_3Sb_2 along the [100] direction (white = Mg_3 , black = Sb_3).

The lanthanoid(III) chloride *cyclo*-tetrasilicates $M_6\text{Cl}_{10}[\text{Si}_4\text{O}_{12}]$ (M=Sm, Gd-Dy): Synthesis, structure and IR investigations

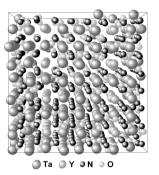
Ingo Hartenbach, Stefan Jagiella and Thomas Schleid Page 2258



Cyclo-oxotetrasilicate unit in $M_6\text{Cl}_{10}[\text{Si}_4\text{O}_{12}]$.

A density-functional and molecular-dynamics study on the physical properties of yttrium-doped tantalum oxvnitride

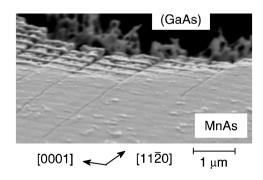
H. Wolff, H. Schilling, M. Lerch and R. Dronskowski *Page 2265*



Structural result of a room-temperature molecular-dynamic simulation of a supercell of $Y_{0.125}Ta_{0.875}O_{0.875}N\square_{0.125}$.

Fabrication of MnAs microstructures on $GaAs(0\,0\,1)$ substrates and their electrical properties

Y. Takagaki, E. Wiebicke, L. Däweritz and K.H. Ploog *Page 2271*

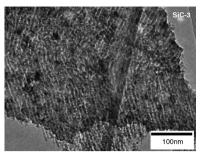


Surface of a partly etched MnAs film grown on a GaAs(001) substrate

Thermal stability of high surface area silicon carbide materials

Piotr Krawiec and Stefan Kaskel

Page 2281

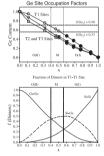


The synthesis of mesoporous silicon carbide by chemical vapor infiltration of dimethyl dichlorosilane into mesoporous silica SBA-15 and subsequent dissolution of the silica matrix was used for the preparation of mesoporous SiC with high specific surface areas up to $830\,\mathrm{m^2\,g^{-1}}$ and pore sizes between 2 and $10\,\mathrm{nm}$.

On the distribution of tetrelide atoms (Si, Ge) in $Gd_5(Si_xGe_{1-x})_4$

Sumohan Misra and Gordon J. Miller

Page 2290

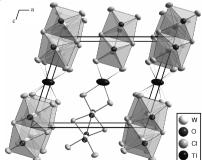


The Ge occupation in each T site in $Gd_5(Si_xGe_{1-x})_4$ is studied as a function of Si concentration, x. The different crystal structures are related to the fractions of Ge–Ge (solid), Si–Ge (dashed) and Si–Si (solid) dimers at the T1–T1 sites.

 $M_{1-x}[W_2O_2X_6]$ with $M=K^+$, Tl^+ , Ag^+ , Hg^{2+} , Pb^{2+} ; X=Cl, Br—A class of mixed valence tungsten (IV,V) compounds with layered structures, W–W bonds and high conductivity

Johannes Beck, Christian Kusterer, Rolf-Dieter Hoffmann and Rainer Pöttgen

Page 2298

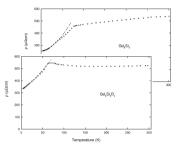


WOCl₃ has a particular reactivity to form non-stoichiometric quaternary compounds $M_{1-x}[W_2O_2Cl_6]$ with mono and divalent cations. The compounds exhibit temperature independent paramagnetism and high electrical conductivity.

Mn_5Si_3 -type host-interstitial boron rare-earth metal silicide compounds RE_5Si_3 : Crystal structures, physical properties and theoretical considerations

Jérome Roger, Mouna Ben Yahia, Volodymyr Babizhetskyy, Joseph Bauer, Stéphane Cordier, Roland Guérin, Kurt Hiebl, Xavier Rocquefelte, Jean-Yves Saillard and Jean-François Halet

Page 2310

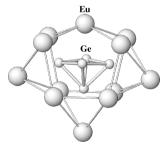


Incorporation of boron in vacant octahedral sites of the Mn₅Si₃-type phase Gd₅Si₃ modifies its resistivity properties.

Barrelane-like germanium clusters in Eu_3Ge_5 : Crystal structure, chemical bonding and physical properties

Sergij Budnyk, Franz Weitzer, Christof Kubata, Yurii Prots, Lev G. Akselrud, Walter Schnelle, Kurt Hiebl, Reinhard Nesper, Frank R. Wagner and Yuri Grin

Page 2329

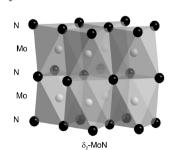


Main building blocks of the crystal structure Eu₃Ge₅ are [Ge₃]⁶⁻ cluster anions surrounded by Eu²⁺ cations. The nearly tetragonal-pyramidal shape is suggested by the interatomic distances. Contrary to that, the bonding analysis with the ELF reveals only two- and three-bonded germanium atoms forming a strongly distorted [111]-barrelane-like cluster.

Synthesis and characterisation of hexagonal molybdenum nitrides

Alexey Yu. Ganin, Lorenz Kienle and Grigori V. Vajenine

Page 2339



Hexagonal δ_2 -MoN with the NiAs-type crystal structure.

Continued

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