MU / S82/12_L March · Vol. 85 · DP17644

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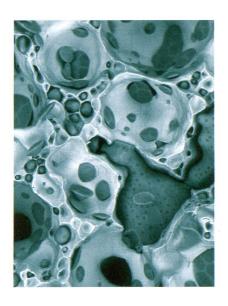












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Cover Photo:

Steel matrix syntactic foams are fascinating materials with great development potential. Their properties include low weight, a metal foam-like response to compressive load and a strength levels exceeding other steel foams. The wide range of matrix/filler combinations possible will allow their fine-tuning to application requirements once structure-property relationships are fully understood. The image shows the fracture surface of a powder metallurgical 316L foam containing 5 wt.-% of glass microspheres, see the article by Dirk Lehmhus and co-workers. Unlike cenospheres, the glass microspheres disintegrate during processing, resulting in a glass phase (medium gray phase, light gray: 316L matrix) embedded in between matrix grains and at inner surface of pores. The latter mark the original position of the microspheres.

Publishing company:

Wiley-VCH Verlag GmbH & Co. KGaA, Boschstraße 12, D-69469 Weinheim, Germany

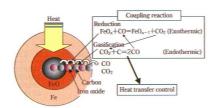
Contents

Review Article

H. M. Ahmed,* N. Viswanathan and B. Bjorkman

Composite Pellets – A Potential Raw Material for Iron-Making

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Using carbon composite pellets as a raw material for iron making processes has the great potential as it can derive the following advantages simultaneously: (i) Replacing expensive coke with relatively less expensive carbon sources, (ii) Producing agglomerates from cheaper raw materials, and (iii) Making the process towards higher carbon utilization. However, they also pose challenges in terms of strength and making the reactions happen at desired conditions. A review on these aspects has been presented here.

Full Paper

D.-Z. Li,* Y.-C. Liu, T.-X. Cui, J.-M. Li, Y.-T. Wang and P.-M. Fu

The Effect of Thermo-Mechanical Processing Parameters on Microstructure and Mechanical Properties of a Low Carbon, High Strength Steel





bainite ferrites BFs and remained austenite films which are located between them and acicular ferrite containing dislocation are presented in figure a. The morphology of martensite M and dislocations D in acicular ferrite are indicated in figure b.

The morphology of carbon-free

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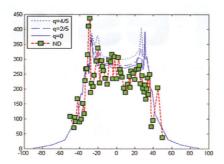
Федеральное государственное бюджетное учреждение науки Центральная научная библиотека Уральского отделения Российской академии наук (ЦНБ УрО РАН)



K. Decroos,* C. Ohms, R. Petrov, M. Seefeldt, F. Verhaeghe and L. Kestens

Influence of Texture on Welding Stress Calculations

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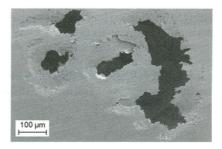


Upon solidification of the fusion zone of an austenitic weld a strong crystallographic texture arises. As a consequence, the bulk fusion zone material becomes transversely isotropic. This paper presents a procedure to define an anisotropic material law for the fusion zone, and to implement that in Finite Element models for welding stress calculations.

H. Pulkkinen,* H. Apajalahti, S. Papula, J. Talonen and H. Hänninen

Pitting Corrosion Resistance of Mn-Alloyed Austenitic Stainless Steels

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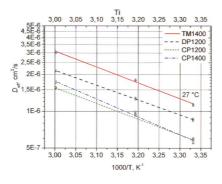


The role of composition, inclusions and precipitates on the pitting corrosion resistance of Mn-alloyed austenitic stainless steels is examined. The Mn-sulfides are observed to act as corrosion pit initiators in chloride containing solutions. As the Cr content is decreased and Ni partially replaced by Mn the pitting corrosion resistance is reduced.

J. Rehrl,* K. Mraczek, A. Pichler and E. Werner

The Impact of Nb, Ti, Zr, B, V, and Mo on the Hydrogen Diffusion in Four Different AHSS/UHSS Microstructures

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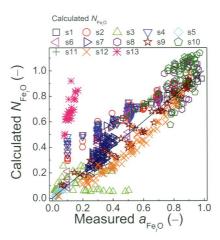
In this work hydrogen diffusion in four non- and microalloyed AHSS/UHSS grades has been investigated. The results indicate that hydrogen diffusion depends significantly on the microstructure adjusted. The tempered martensitic grade shows the highest diffusivity, followed by the dual-phase steel. Complexphase steel grades reveal the lowest diffusivity. However, precipitates and substitutional atoms of microalloying elements seem to have a negligible effect on hydrogen diffusion.



X.-M. Yang,* M. Zhang, J.-L. Zhang, P.-C. Li, J.-Y. Li and J. Zhang

Representation of Oxidation Ability for Metallurgical Slags Based on the Ion and Molecule Coexistence Theory

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To calculate the defined comprehensive mass action concentration $N_{\rm Fe,O}$ of iron oxides in the selected slag systems, a thermodynamic model for calculating the mass action concentrations of structural units or ion couples in CaO-SiO₂-MgO-FeO-Fe₂O₂-MnO- $Al_2O_3-P_2O_5$ type slags, i.e., the IMCT- N_1 model, has been developed. The cover shows the reliable agreement between the calculated comprehensive mass action concentration $N_{\text{Fe},0}$ of iron oxides by the developed $\dot{I}MCT-N_i$ model and reported activity a_{Fe,O} of iron oxides in 13 slag systems at different temperatures. The determined comprehensive mass action concentration $N_{\rm Fe,O}$ of iron oxides can be successfully applied to predict the oxidation ability of metallurgical slags.

A. Tilliander,* L. T. I. Jonsson and P. G. Jönsson

A Three-Dimensional Three-Phase Model of Gas Injection in AOD Converters

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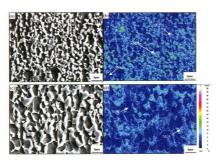


A 3D 3-phase (steel, slag and gas) AOD model with an industrial relevant geometry including six nozzles has been used to predict fluid flow, turbulence and bubble characteristics as well as fluid-slag dispersion. Also, two different gas flow rates have been simulated. This new findings opens up for further investigations of gas-metal reactions in the AOD converter.

C. Wang, H. Ding,* Z. Tang, J. Zhang and H. Di

The Effect of Intercritical Annealing on the Microstructure and Mechanical Properties in a High Strength TRIP Steel with C-Partitioning

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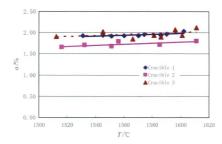


In this study, the effect of intercritical annealing on the microstructure and mechanical properties with C-partitioning is investigated for a high strength transformation-induced plasticity steel (TRIP steel). The partitioning of C to austenite during intercritical annealing are the main contributions to the austenite stability is confirmed by direct experimental evidence (WDS). After heat treatment, C atoms mainly exist among the granular retained austenite and the M/A islands.



L. Zhong,* M. Zeze and K. Mukai Density and Molar Volume of Liquid Iron Containing Nb

395

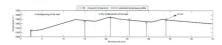


The linear expansion of the alumina crucibles used in this density measurement is about 1.65–2.07 at the temperature range from 1510 to 1610°C. The linear expansion increases linearly with an increase in temperature. Crucible 1 and 3 seem to have the same expansion coefficients, but crucible 2 has a lower expansion coefficient than those of crucible 1 and 3.

W. Lv, Z. Mao,* P. Yuan and M. Jia

Pruned Bagging Aggregated Hybrid Prediction Models for Forecasting the Steel Temperature in Ladle Furnace

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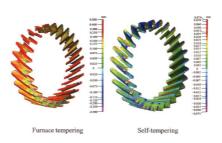


A novel prediction model is developed for real-time forecasting the steel temperature during LF refining process. In order to obtain good continuous prediction capability, a hybrid predictor is firstly developed by combining first principle method with data-driven modeling technique. Thereafter, a modified ensemble learning algorithm is introduced for further boosting prediction accuracy.

D. Rodman,* F. Nürnberger, A. Dalinger, M. Schaper, C. Krause, M. Kästner and E. Reithmeier

Tempering Induction Hardened 42CrMo4 Steel Helical Gearwheels from Residual Heat Using Spray Cooling

_ 415

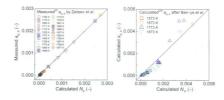


During induction hardening of helical gearwheels made of 42CrMo4 heat-treating steel using spray cooling, a self-tempering was performed. The result of the heat treatment is analyzed with the aid of hardness, residual stress, and distortion measurements. The obtained hardening results demonstrate the equivalence of self-tempering and conventional furnace tempering.

X.-M. Yang,* P.-C. Li, J.-Y. Li, M. Zhang, J.-L. Zhang and J. Zhang

Representation Reaction Abilities of Structural Units and Related Thermodynamic Properties in Fe–P Binary Melts Based on the Atom-Molecule Coexistence Theory

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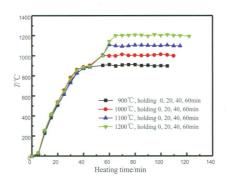
A thermodynamic model for calculating the mass action concentrations of structural units in Fe-P binary melts based on the atom-molecule coexistence theory, i.e., AMCT-N. model, has been developed and verified through comparing with the reported activities of both P and Fe in Fe-P binary melts with mole fraction x_0 of P < 0.33 in a temperature from 1406 K to 1973 K. The cover shows that the calculated mass action concentration $N_{\rm p}$ of P can correlate a very good 1:1 corresponding relationship with the reported activity a_{RP} of P relative to pure liquid P(l) as standard state.



J.-j. Hao, J. Chen,* P.-d. Han, K. Li and H.-f. Sun

Solid Phase Decarburization Kinetics of High-Carbon Ferrochrome Powders in the Microwave Field

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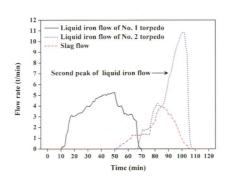


This paper deals with the kinetics of solid phase decarburization of HCFPs in microwave heating filed. Apparent activation energy of the solid phase decarburization reaction is calculated and restrictive step and the mechanism is also obtained. The optimum conditions of solid phase decarburization of HCFP by microwave heating are also concluded.

K. Du,* S. Wu, M. Kou, W. Shen and Z. Zhang

Numerical Simulation of Liquid Iron Flow and Heat Transfer in the Hearth of COREX Melter Gasifier during Tapping Process

__ 466



In comparison with blast furnace, the second peak of liquid iron flow at the later stage of tapping, which mainly results from the expansion of taphole diameter, is a distinguishing feature of COREX melter gasifier hearth. The practical experience shows that depressurization tapping and improving the clay properties are effective measures to minimize the second peak of liquid iron flow.

A. Jahn, K.-P. Steinhoff, T. Dubberstein,* P. Franke, M. Weider, S. Wolf, A. Kovalev, A. Glage, A. Weiß, W. Schärfl, K. Eigenfeld, L. Krüger and P. R. Scheller

Phosphor Alloyed Cr–Mn–Ni Austenitic As-cast Stainless Steel with TRIP/ TWIP Effect

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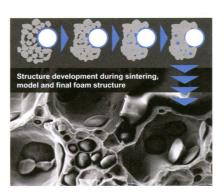


The influence of increasing phosphor content on the flowability, the effect on the infiltration of ZrO_2 open foam ceramics, the martensite formation as well as the resulting mechanical properties and the surface tension of high alloyed Cr–Mn–Ni as-cast steels are studied. The elongation properties are not influenced by increasing phosphor content.

J. Weise, D. Lehmhus,* J. Baumeister, R. Kun, M. Bayoumi and M. Busse

Production and Properties of 316L Stainless Steel Cellular Materials and Syntactic Foams

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316L steel matrix syntactic foams are produced via MIM and using soda-lime borosilicate hollow glass microspheres as filler. Structural development is evaluated using metallographic, image analysis and FIB preparation as well as EDX analysis techniques. Mechanical behavior is studied under quasistatic compressive and tensile load and qualitatively linked to structural features.