PAPER • OPEN ACCESS

High-temperature Mechanical Properties and Microstructure of ZrTiHfNbMo_x (x=0.5, 1.0, 1.5) Refractory High Entropy Alloys

To cite this article: Y W Chen et al 2018 IOP Conf. Ser.: Mater. Sci. Eng. 359 012033

View the article online for updates and enhancements.

You may also like

- First principles calculations to investigate Li-based quaternary Heusler compounds LiHfCoX (X = Ge, Sn) for thermoelectric applications

Tavneet Kaur, Jaspal Singh, Megha Goyal et al.

- <u>Revealing photoluminescence and</u> nonlinear optical absorption characteristics of PbMo_{0.75}W_{0.25}O₄ single crystal for optical limiting applications Anl Doan, Ahmet Karatay, Mehmet Isik et al.
- First-principles study of the phase stability and elastic properties of TiVNbMoM (M = Al, Sc, Ni and Cu) high entropy alloys N Al-Zoubi





DISCOVER how sustainability intersects with electrochemistry & solid state science research



This content was downloaded from IP address 3.143.168.172 on 28/04/2024 at 19:33

High-temperature Mechanical Properties and Microstructure of ZrTiHfNbMo_x (x=0.5, 1.0, 1.5) Refractory High Entropy Alloys

Y W Chen, Y K Li^{*}, X W Cheng, C Wu and B Cheng

School of Materials Science and Engineering, Beijing Institute of Technology, Beijing 100081, China

National Key Laboratory of Science and Technology on Materials under Shock and Impact

E-mails: 2861209902@qq.com; liyunkai@bit.edu.cn

Abstract: Refractory high entropy alloys (RHEAs), with excellent properties at high temperature, have several applications. In this work, the ZrTiHfNbMo_x (x=0.5, 1.0, 1.5) alloys were prepared by arc melting. All these alloys form body centered cubic (BCC) structure without other intermediate phases. The Mo element contributes to the strength of alloys at high temperature, but too much of Mo decreases the plasticity severely and enhances the strength. The ZrTiHfNbMo alloy, whose compressive stress is 1099 MPa at 800° C, is a promising material for high-temperature applications.

1.Introduction

In recent years, RHEAs were proposed due to their extraordinary properties in the high temperature application field[1-3]. Contrasted with the traditional superalloys,RHEAs containmore than five elements with near equal mole ratios (5%-35%)and form the single microstructure [4, 5]. Because the elements of RHEAsare in the group IV, V and VI with the melting point over 1650 °C, the operating-temperature range of these alloysare more than Al-Ni superalloy [6, 7]. The special performances of RHEAs including high configural entropy, sluggish diffusion, the severe lattice distortionand unique cocktail effect [8-11] contribute to the high strength even in the extreme conditions, which become promising candidates in high temperature materials field.

2. Experimental

The ZrTiHfNbMo_x(x=0.5, 1.0, 2.0) alloys were prepared by arc melting under the purified argon atmosphere, and the cooling system employed the water-cooled copper crucible. The as-cast crystalline structure was investigated by X-ray diffractometer (XRD, D8 advance) with K_{α} radiation. The high temperature compressive performances were performed on a Gleeble-3500 thermal simulator, each alloy was heated 800°C and 900°C in turns, the heating rate was setting as 300°C/min and the holding time was 3 minutes. A scanning electron microscope (SEM, HITACHIS4800) was adopted to analyze the microstructure with the secondary electron image.

3. Results and discussion

3.1. Crystal structure and microstructure analysis

The X-ray diffraction patterns of these as-cast ZrTiHfNbMo_xalloys are shown in the figure 1.According to the analysis through JADE 5.0 software, all the major peaks of these alloys are consistent with the single phase of body center cubic without any other precipitated phase. The diffraction peaks aresharp, which indicates that these alloys have high crystallinity. With the increase of Mo content (the radius of Mo element is apparently smaller than other element), the diffraction peak shift to the low theta due to the crystal lattice become smaller.



Figure 1. X-ray diffraction patterns of the RHEAs

3.2. High temperature compression performance analysis

The compressive stress-strain curves of ZrTiHfNbMo_x alloys at high temperature are displayed in the figure 2. The process of stress follows that increase rapidly, reduction and remaining a stable value with the deformation crease. The high-temperature stress of the ZrTiHfNbMo_{0.5} alloy is relatively poor, which is 585 MPa at 800°C and 288 MPa at 900°C. When adjusting the content of Mo element, the stress of ZrTiHfNbMo alloy is1099MPa at 800°Cand804MPaat 900°C. The Mo element has the highest melting point and strength among these elements, which contributes to the properties of softening resistance. As the Mo content continues to increase, the stress of ZrTiHfNbMo_{2.0} alloy is 1155MPa at 800°C, and 897MPa at 900°C, but the strain is less than 15% at 800°C contrasted with

thestrain of ZrTiHfNbMo_{0.5} and ZrTiHfNbMo alloy more than 40%, respectively. The excessive Mo element made the lattice distortion increase and the dislocation slides difficultly even at high temperature. As the high temperature could provide much more energy, then it would promote the movement of dislocation, vacancy and grain boundary, which could decrease the stress.

The compression scanning electron microscopies of ZrTiHfNbMo_xalloys at high temperature are shown in the figure 3. Compared the figure.3 (a) and (b) with thefigure.3 (d) and (e), the size of grains become larger and the intercrystalline crack is more serious as the temperature increases, part of the dendrites were fused. The fracture feature of ZrTiHfNbMo_{2.0} alloys is the river-like appearance in the figure.3(c), which reveals that the poor plasticity at 800°C.



Figure 2. High-temperature compressive stress-strain curves of these alloys: (a) Mo_{0.5}; (b) Mo_{1.0}; (c) Mo_{2.0}



Figure 3.The high-temperature fracture scanning electron microscopies of ZrTiHfNbMo_xalloys. Mo_{0.5}: (a) at 800°C and (d) at 900°C; Mo_{1.0}: (b) at 800°C and (e) at 900°C; Mo_{2.0}: (c) at 800°C and (f) at 900°C;

4. Conclusion

In this paper, the ZrTiHfNbMo_x (x=0.5, 1.0, 2.0) refractoryHEAs were melted and the high-temperature properties were analyzed. These alloys formed the body centered cubic structure without other phases. The high temperature stressof these alloys is enhanced with the Mo element increase (Mo content from 0.5 to 1.0). The strength of ZrTiHfNbMo_{2.0} enhances a few and their plasticity decreases rapidlyas the Mo element continues to increase, and the fracture displays the river-like appearance at 800 °C. The reasonable content of Mo element is contributed to the high temperature mechanical properties, and the HEAs with balance between strength and plasticityat high temperature can be acquired through controlling the Mo content addition.

References

- C.-Y. Hsu, J.-W. Yeh, S.-K. Chen, T.-T. Shun, Wear resistance and hightemperature compression strength of FCC CuCoNiCrAl0.5Fe alloy with boron addition, Metall. Mater. Trans. A 35A (2004) 1465-1469.
- [2] J.-W. Yeh, S.-K. Chen, J.-W. Gan, S.-J. Lin, T.-S. Chin, T.-T. Shun, C.-H. Tsau, S.Y. Chang, Formation of simple crystal structures in Cu-Co-Ni-Cr-Al-Fe-Ti-V alloys with multiprincipal metallic elements, Metall. Mater. Trans. A 35A (2004) 2533-2536
- [3] J.-W. Yeh, S.-K. Chen, S.-J. Lin, J.-Y. Gan, T.-S. Chin, T.-T. Shun, C.-H. Tsau, S.Y. Chang, Nanostructured high-entropy alloys with multiple principal elements: Novel alloy design concepts and outcomes, Adv. Eng. Mater. 6 (2004) 299-303.
- [4] Yeh, J.-W., et al., Anomalous decrease in X-ray diffraction intensities of Cu-Ni-Al-Co-Cr-Fe-Si alloy systems with multi-principal elements. Materials Chemistry and Physics, 2007. **103**(1): p. 41-46.
- [5] Cantor, B., et al., *Microstructural development in equiatomic multicomponent alloys*. Materials Science and Engineering: A, 2004. **375-377**: p. 213-218.
- [6] Senkov, O.N., et al., *Development of a Refractory High Entropy Superalloy*. Entropy, 2016. **18**(3).
- [7] Senkov, O.N., et al., *Microstructure and room temperature properties of a high-entropy TaNbHfZrTi alloy.* Journal of Alloys and Compounds, 2011. **509**(20): p. 6043-6048.
- [8] Chuang, M.H., et al., *Microstructure and wear behavior of AlxCo1.5CrFeNi1.5Tiy high-entropy alloys.* Acta Materialia, 2011. **59**(16): p. 6308-6317.
- [9] Zhou, Y.J., et al., *Solid solution alloys of AlCoCrFeNiTix with excellent room-temperature mechanical properties.* Applied Physics Letters, 2007. **90**(18): p. 181904.
- [10] Yang, X. and Y. Zhang, *Prediction of high-entropy stabilized solid-solution in multi-component alloys.* Materials Chemistry and Physics, 2012. **132**(2-3): p. 233-238.
- [11] JianBin Guo, YunKai Li, *Study on microstructure and mechanical properties of five-element refractory high entropy alloys*, D. Beijing Institute of Technology, 2015.