

## L Edge Absorption Measurements of $\text{ZrO}_2$ and $\text{Y}_2\text{O}_3$ at BL13

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$L_{2,3}$  edges of Y and Zr are known to have sharp structures depending on chemical composition [1, 2]. We have performed measurements of not only the  $L_{2,3}$  edges but also the  $L_1$  edge of  $\text{Y}_2\text{O}_3$  and  $\text{ZrO}_2$  in which Y and Zr are formally in  $\text{Y}^{3+}$  and  $\text{Zr}^{4+}$  states, both having the same electronic configuration as Kr.

Measurements have been carried out at BL-13 of the SR Center of Ritsumeikan University in the total electron yield method. Powdered samples of  $\text{Y}_2\text{O}_3$  (Kojundo Chemical Lab., 99.9 %) and  $\text{ZrO}_2$  (NewMet, 99.8 %) were attached on the carbon tape.

In Fig. 1,  $L_{2,3}$  spectra of  $\text{Y}_2\text{O}_3$  and  $\text{ZrO}_2$  are shown. The  $L_3$  lineshape of  $\text{Y}_2\text{O}_3$  is similar with the reported one [1]. The  $L_{2,3}$  lineshape of  $\text{ZrO}_2$  is similar with the reported spectrum of monoclinic  $\text{ZrO}_2$  (m- $\text{ZrO}_2$ ), which is the room-temperature phase [2]. In each of the  $L_3$  and  $L_2$  edges, two structures roughly 2 eV apart, are seen. The intensity of the higher energy structure is relatively stronger at the  $L_3$  edge than at the  $L_2$  edge. Furthermore, the intensity of the higher energy structure is relatively stronger for  $\text{ZrO}_2$  than for  $\text{Y}_2\text{O}_3$ .

According to Ikeno et al.[2], the above mentioned structures in the  $L_{2,3}$  edges mainly reflect the ligand-field splitting of the unoccupied 4d state of Y and Zr, whereas multiplet effects play secondary roles. The 2-eV split lineshape of Y- $L_{2,3}$  edges in  $\text{Y}_2\text{O}_3$ , which resembles the lineshape of the reported Zr- $L_{2,3}$  edges in  $\text{SrZrO}_3$ [2], reflects the  $t_{2g}$ - $e_g$  splitting because of the octahedral (six-fold) coordination [3]. On the other hand, the lineshape of m- $\text{ZrO}_2$  was explained by splitting into five levels due to seven-fold coordination with lower symmetry [2].

Y and Zr  $L_1$  spectra of  $\text{Y}_2\text{O}_3$  and  $\text{ZrO}_2$  are shown in Fig. 2. In the Y  $L_1$  spectrum, a structure is found at about 10 eV above the main peak. In the Zr  $L_1$  spectrum a broader and weaker hump seems to exist around 10 eV above the main peak. These lineshapes probably reflect the unoccupied Y and Zr 5p states.

### References

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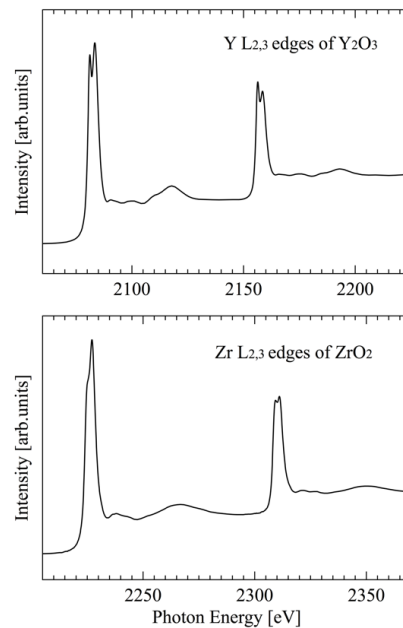


Fig. 1. Total electron yield spectra at Y and Zr  $L_{2,3}$  edges of  $\text{Y}_2\text{O}_3$  and  $\text{ZrO}_2$ .

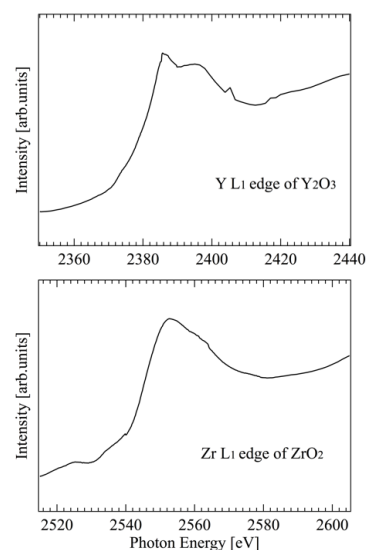


Fig. 2. Total electron yield spectra at Y and Zr  $L_1$  edge of  $\text{Y}_2\text{O}_3$  and  $\text{ZrO}_2$ .