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9
Apêndice

9.1. Difractogramas indicando o respectivo ajuste Rietveld.

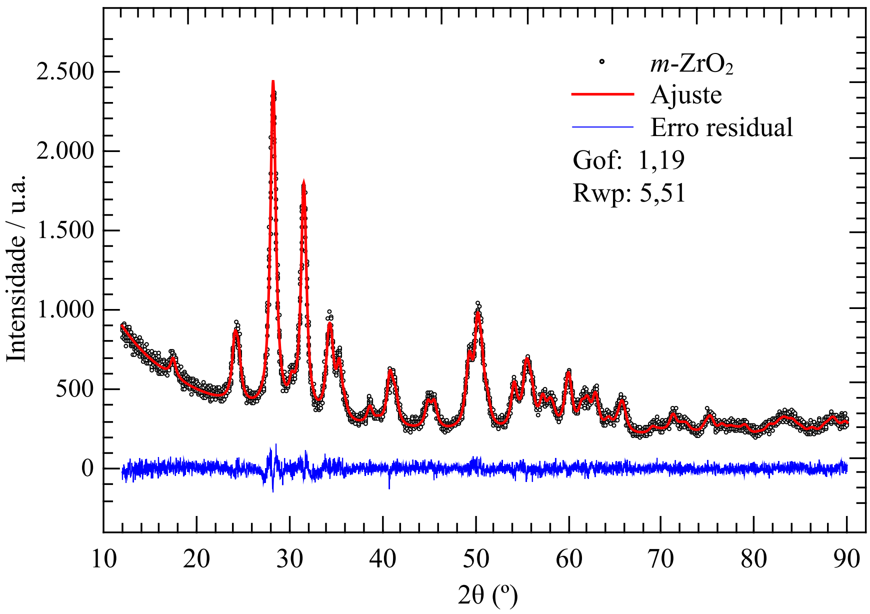

Figura A1. Difractograma da amostra *m*-ZrO_2 e ajuste pelo método de *Rietveld*

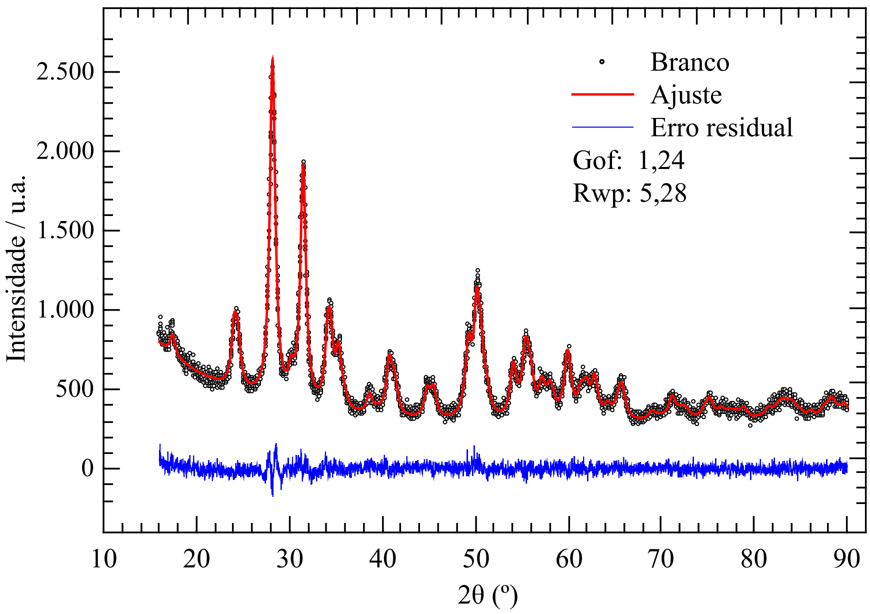

Figura A2. Difractograma da amostra *branco* e ajuste da fase *m*-ZrO_2 pelo método de *Rietveld*

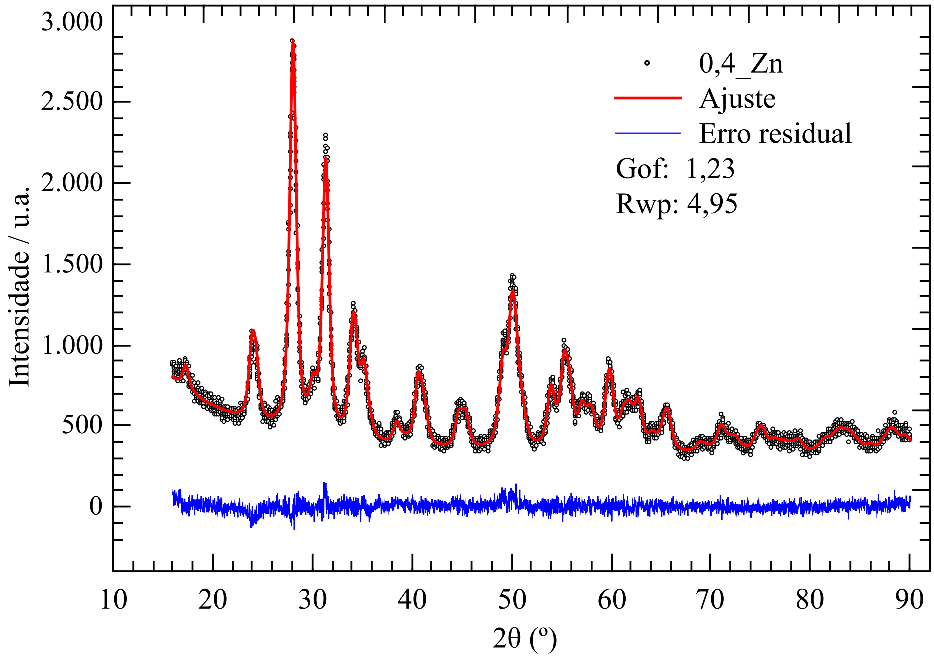

Figura A3. Difractograma da amostra 0,4_Zn e ajuste da fase *m*-ZrO_2 pelo método de *Rietveld*

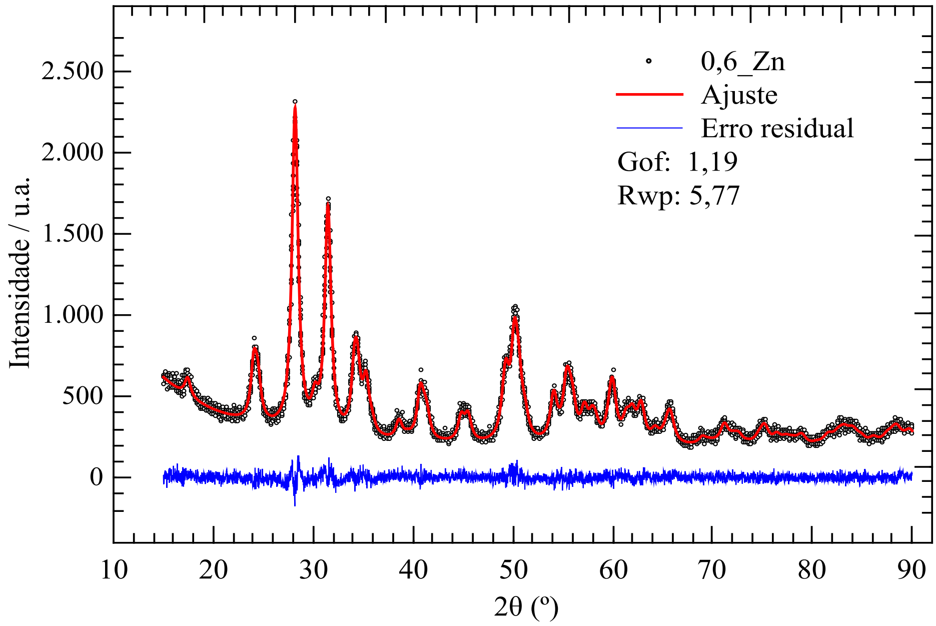

Figura A4. Difractograma da amostra 0,6_Zn e ajuste da fase *m*-ZrO_2 pelo método de *Rietveld*

9.2. Espectros XPS indicando valores de energia de ligação do zinco

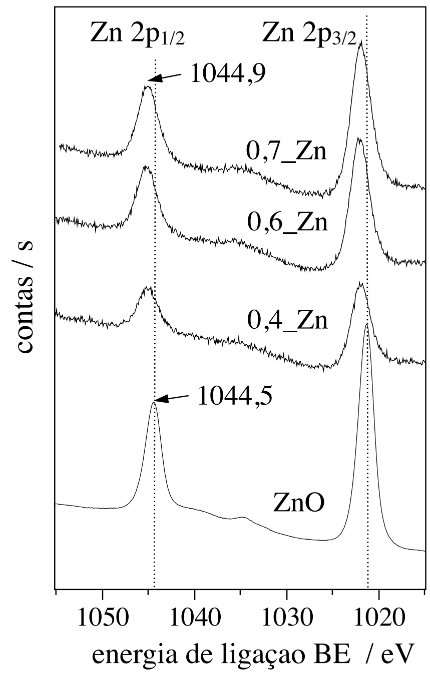

Figura A5. Energias de ligação Zn 2p $_{1/2}$ e Zn 2p $_{3/2}$ dos óxidos mistos sintetizados e de uma amostra comercial de ZnO.

9.3. Deconvolução dos espectros de TPD-CO_2.

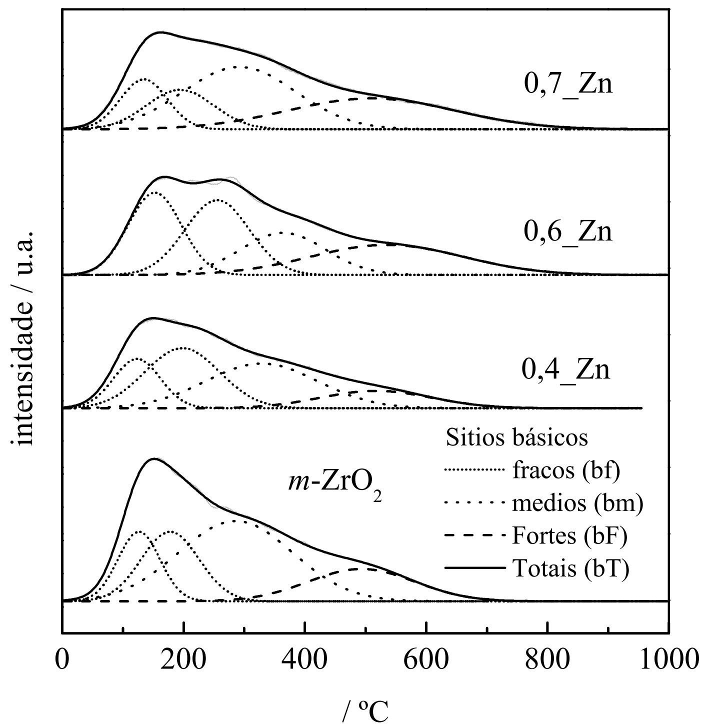

Figura A6. Deconvolução dos espectros de TPD-CO_2 indicando densidade de sítios básicos.

9.4. Deconvolução dos espectros de TPD-NH_3.

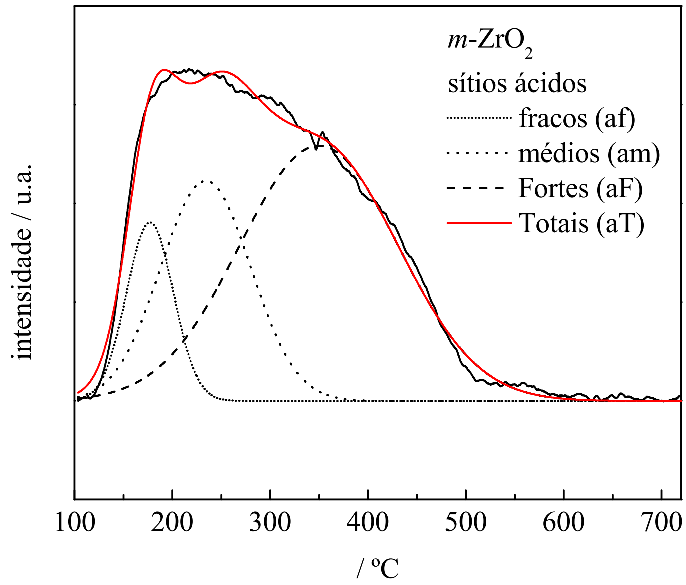

Figura A7. Deconvolução dos espectros de TPD-NH_3 realizado na amostra *m*-ZrO_2, indicando densidade de sítios ácidos.

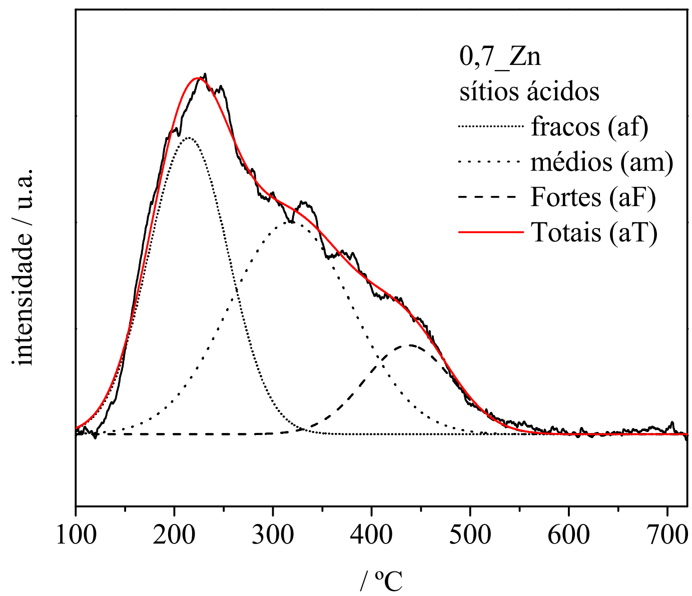

Figura A8. Deconvolução dos espectros de TPD-NH_3 realizado no óxido misto 0,7_Zn, indicando densidade de sítios ácidos.