Multiscale phenomena in molecular matter



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Optimalization of anodization parameters for nanoporous Al₂O₃ templates fabrication

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Nanoporous Al_2O_3 have a wide range of applications as a sensors, biosensors or energy storage devices. These templates are attractive due to their low fabrication costs.

The Al oxides with hexagonal close packed porous structure can be prepared by two step anodizing process with features tunable by the electrolyte composition and voltage. It was already shown [1], [2] that degree of pore ordering depends on first anodization time and the initial surface roughness. This work is focused on Al foil surface pretreatment procedure and first-step anodization parameters influence on order parameter of pore structure.

A series of porous anodic aluminium oxide samples were prepared by two-step

anodization of the high purity (99.999%) a luminium foil in oxalic acid at temperature \sim 17 °C. The preparation of Al foils consisted of annealing at 600 °C for 48 hours, oxide layer removal and fine chemical polishing. It was found that such a sequence allows to obtain a surface with roughness of approximately 10 nm.

The time of the first anodization was varied from 10 min to 15 hours, while the second anodization time was 10 minutes. Pores ordering was estimated by self-ordering parameter determined from the number of violations in hexagonal symmetry for 100 pores, and by comparison of Fast Fourier Transforms (FFT) obtained for SEM images of sample surfaces. It was found that increasing of first-step anodization time from 10 minutes to 2.5 hours

significantly enhances self-ordering parameter from 0.66 to 0.87, respectively. Improvement of FFTs'symmetry confirms ordering enhancement. Further extension of anodization time do not affect the ordering. Therefore, the time of 2.5 h for first-step anodization followed by the annealing and chemical polishing of Al foil was found to be an optimum for well-ordered templates fabrication.

References

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