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Physical and Electrochemical Properties  
of  $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$  Obtained From Different  
Preparation Method

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**Physical and Electrochemical Properties of  $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$   
Obtained From Different Preparation Method**

삼성전관 기술본부

에너지LAB

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Energy Lab

## Objective

To Optimize Preparation Method and to Investigate Cycling Behavior of the  $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$  ( $x = 0.1, 0.2$  and  $0.3$ ) Cathode Materials

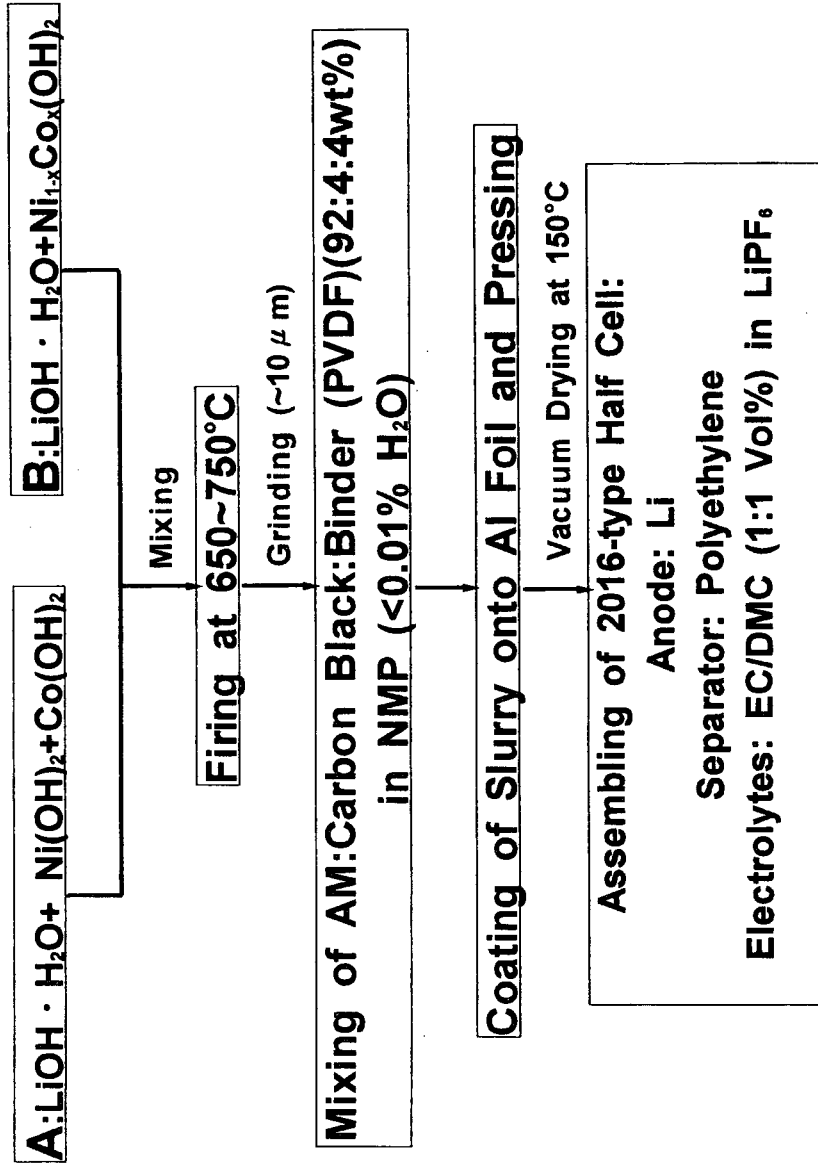
### 1. Optimization of Synthesis Method

- No Formation of  $\text{CO}_2$  Gas
- Low Cost
- Mass Producible: Increasing Degree of Ni and Co Mixing

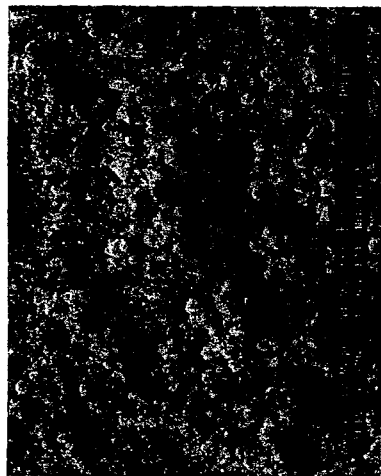
### 2. Investigation of Cycling Behavior of $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$

- High Specific Capacity
- Good Capacity Retention at 1C ( $=180\text{mA/g}$ )
- Low Irreversible Capacity ( $<20\text{mAh/g}$ )

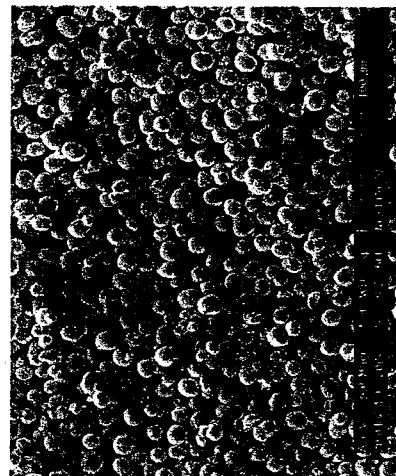
# Experimental



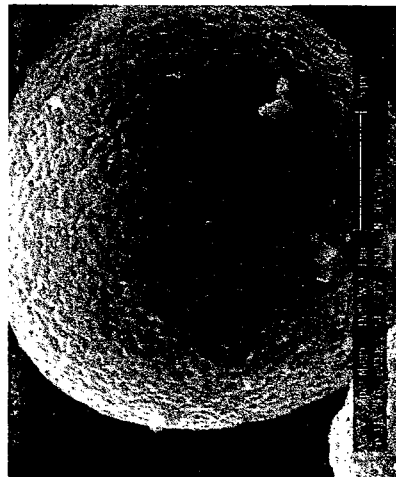
**SEM Pictures of  $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$  Prepared by Different Methods**



**Method A**  
Tap density:  
2.6cc/g

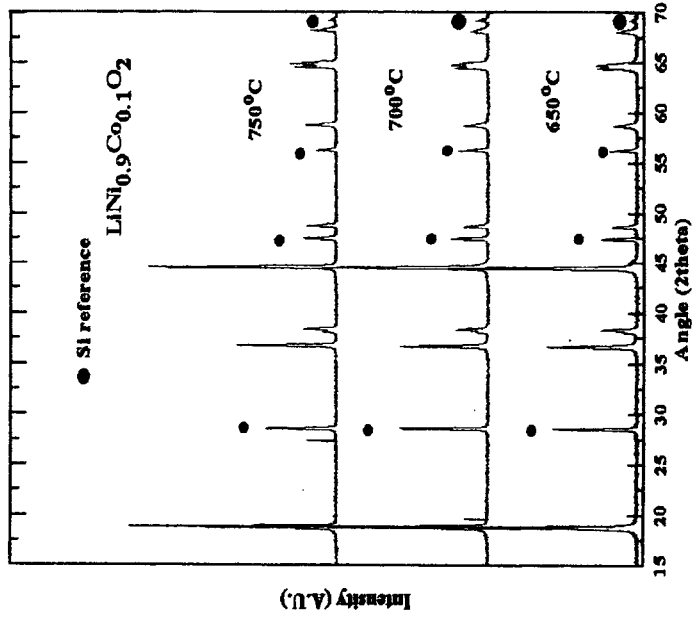


**Method B**  
Tap density:  
2.9cc/g

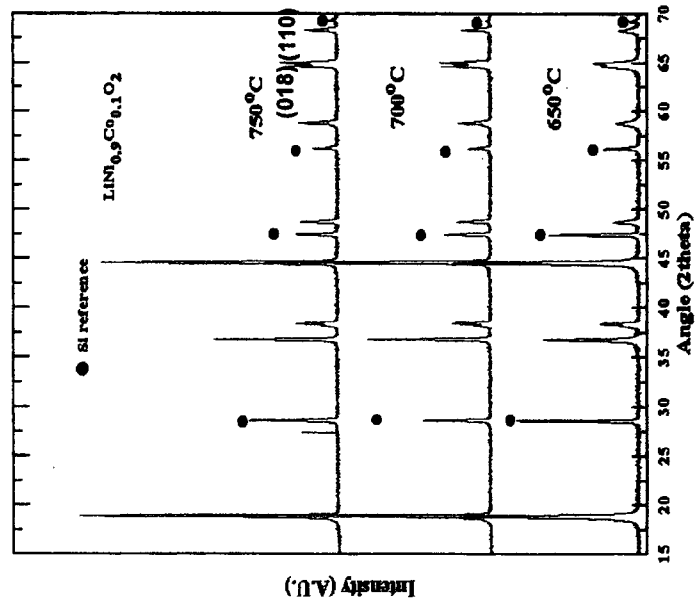


### Comparison of the Powder XRD Patterns

**Method A**

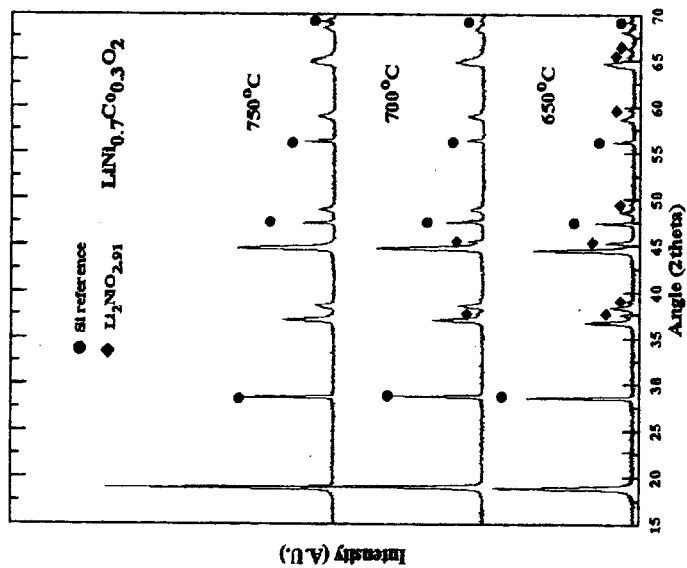


**Method B**

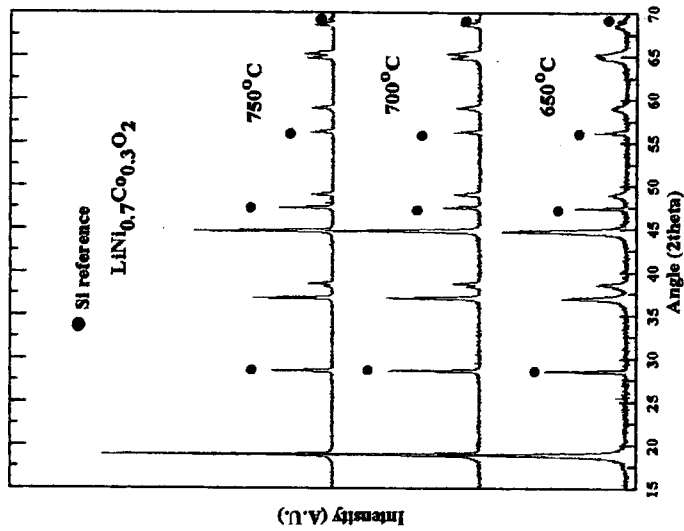


## Comparison of the Powder XRD Patterns

**Method A**



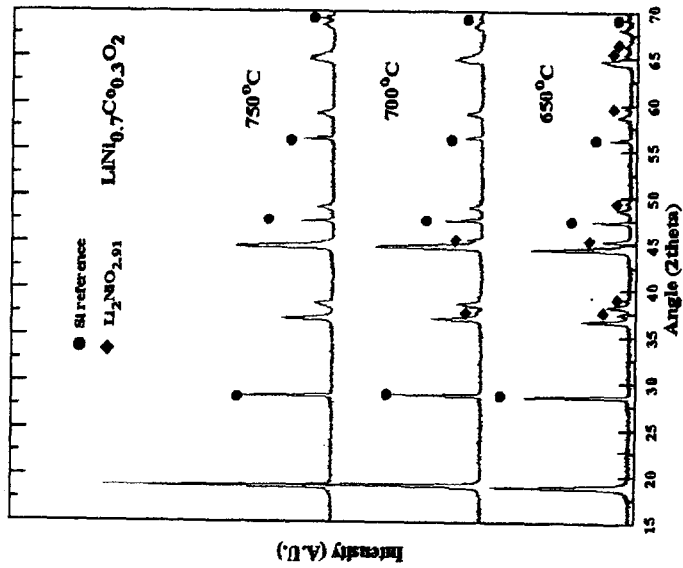
**Method B**



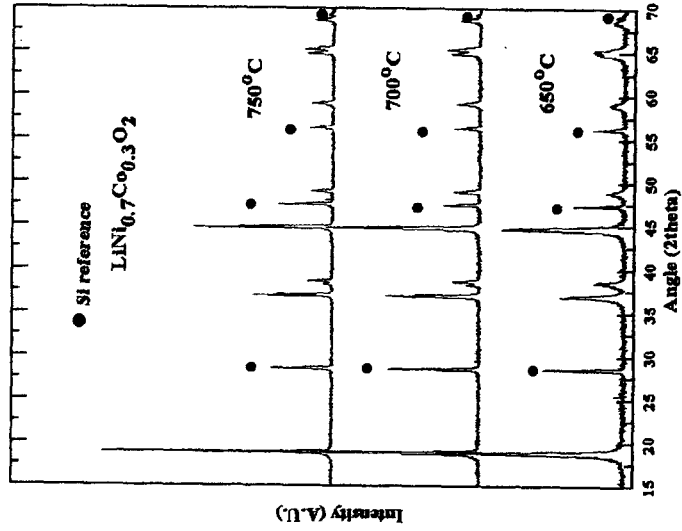


## Comparison of the Powder XRD Patterns

**Method A**



**Method B**



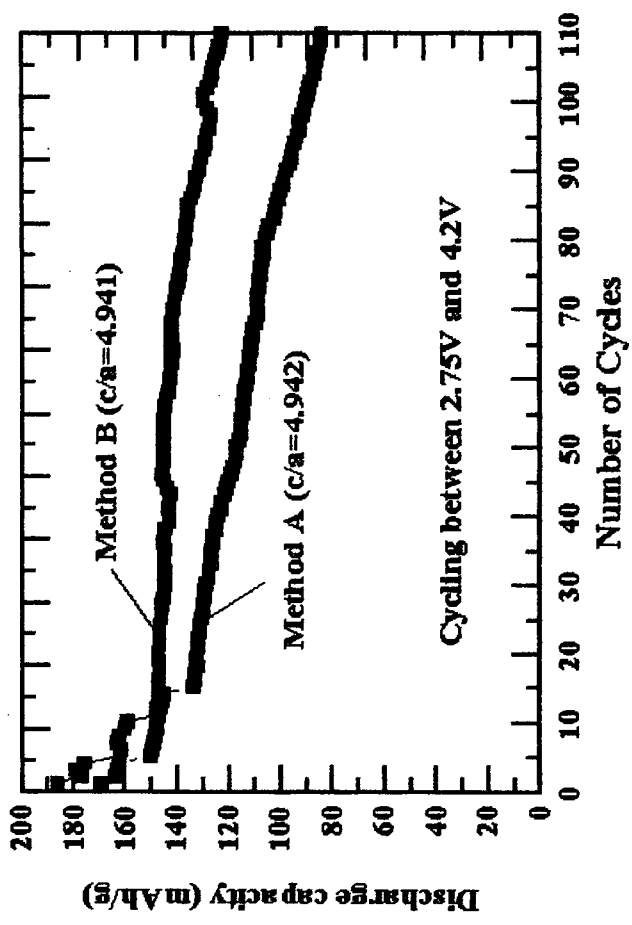
## Comparison of $a/c$ Values in the $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$ Materials Depending on Firing Temperature

Method	A			B		
	$\text{LiNi}_{0.9}\text{Co}_{0.1}\text{O}_2$	$\text{LiNi}_{0.8}\text{Co}_{0.2}\text{O}_2$	$\text{LiNi}_{0.7}\text{Co}_{0.3}\text{O}_2$	$\text{LiNi}_{0.9}\text{Co}_{0.1}\text{O}_2$	$\text{LiNi}_{0.8}\text{Co}_{0.2}\text{O}_2$	$\text{LiNi}_{0.7}\text{Co}_{0.3}\text{O}_2$
650°C	4.930	4.918	4.909	4.933	4.934	4.938
700°C	4.942		4.924	4.936	4.94	4.951
750°C	4.942		4.910	4.934	4.941 <sup>#</sup>	4.948

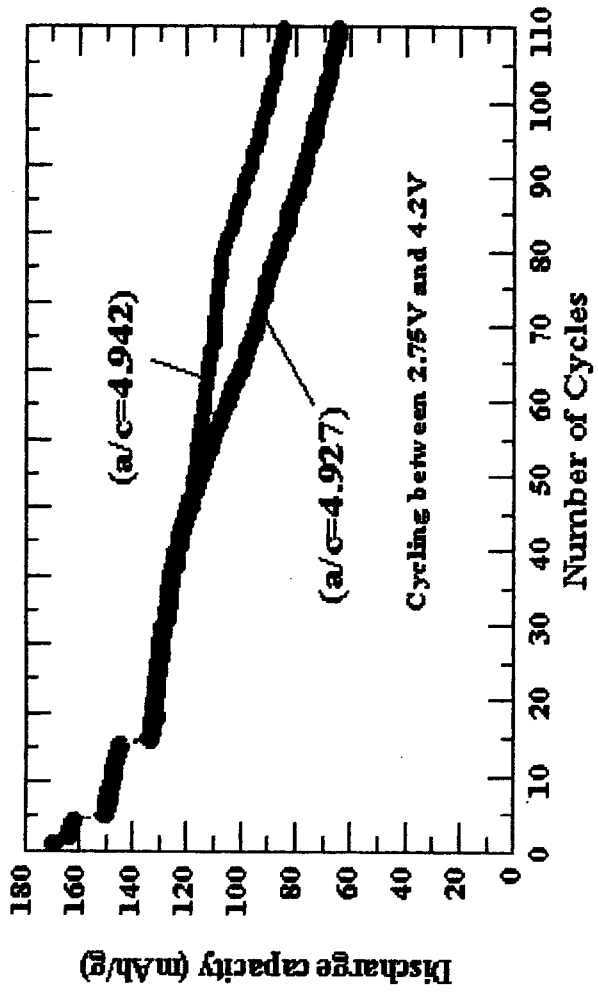
\*  $\text{LiCoO}_2$ : 4.99,  $\text{LiNiO}_2$ : 4.935

\*  $c/a$  is an indicative of structure anisotropy (cubic lattice:  $a/c=4.899$ )

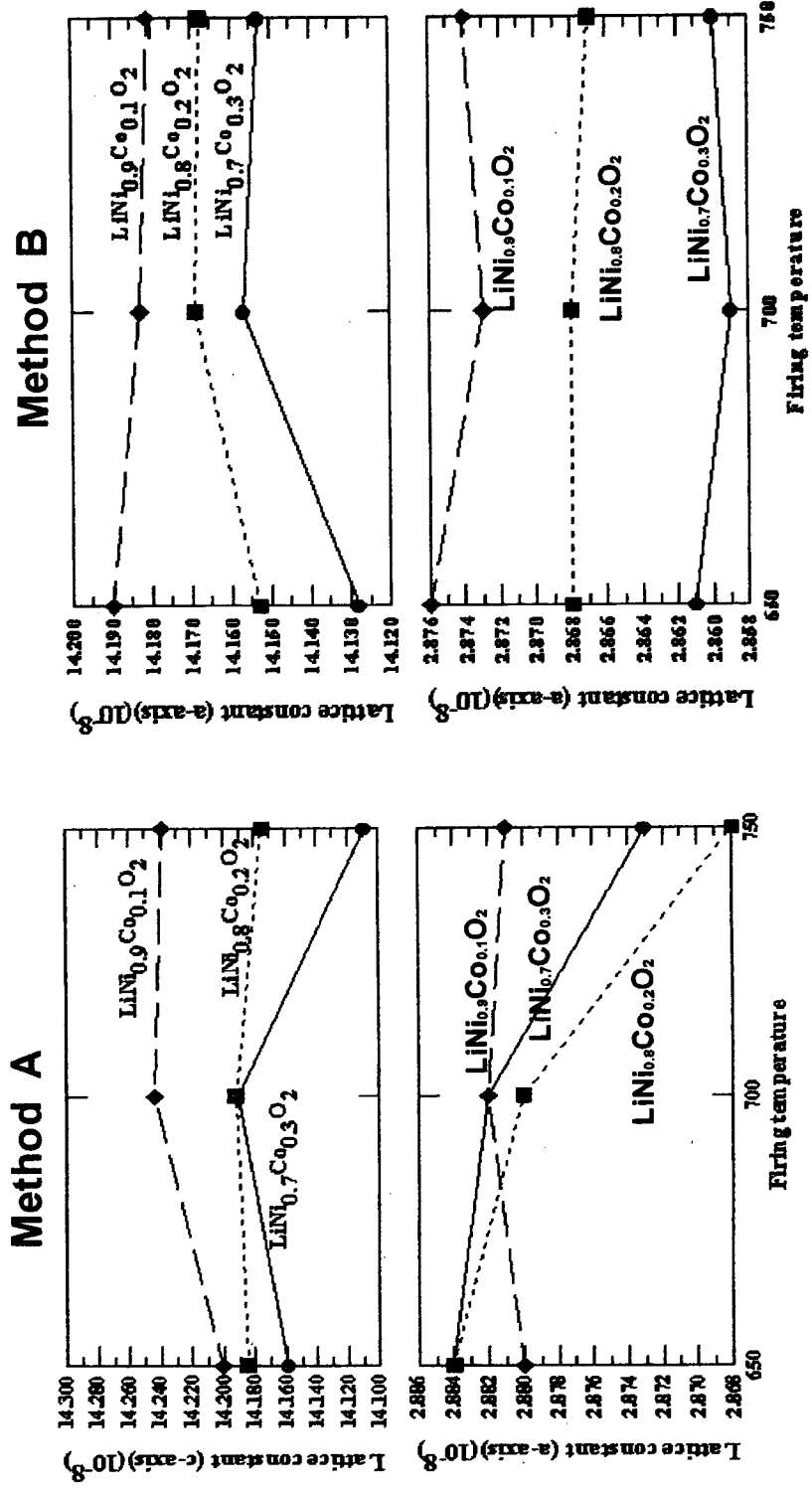
### Cycling Behavior of $\text{LiNi}_{0.8}\text{Co}_{0.2}\text{O}_2$ Prepared by Different Methods



### Cycling Behavior of $\text{LiNi}_{0.8}\text{Co}_{0.2}\text{O}_2$ (A Methode) at Different C-rates

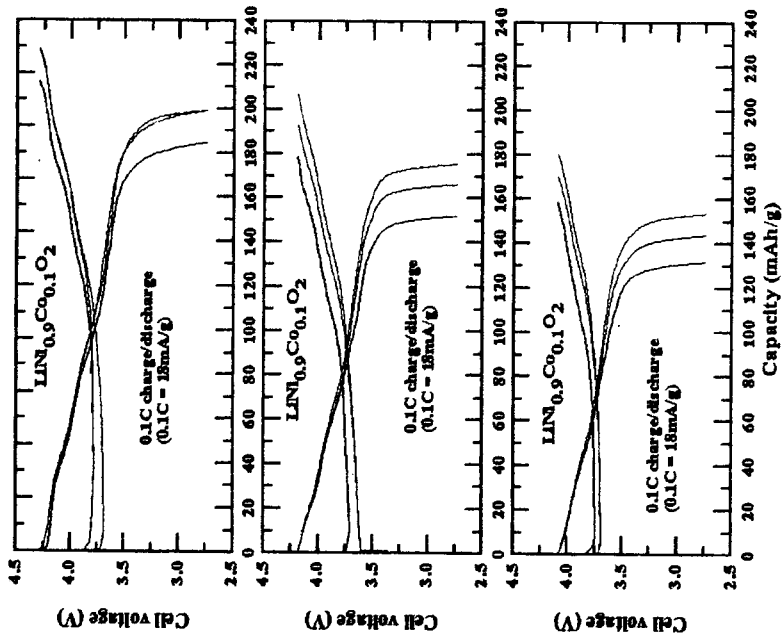


### Comparison of Lattice Parameters as a Function of Firing Temperature

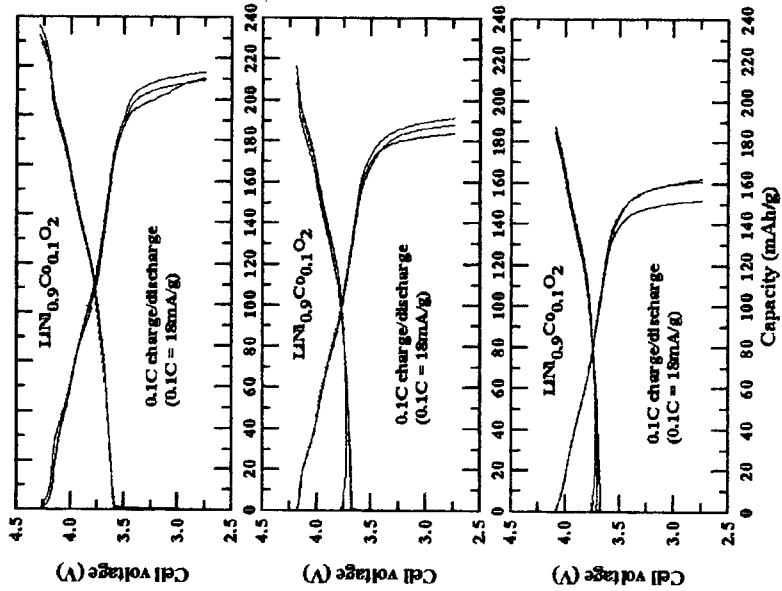


**Comparison of the Cycling Curves in the First Cycle at Different Cut-off Voltages**

**Method A**

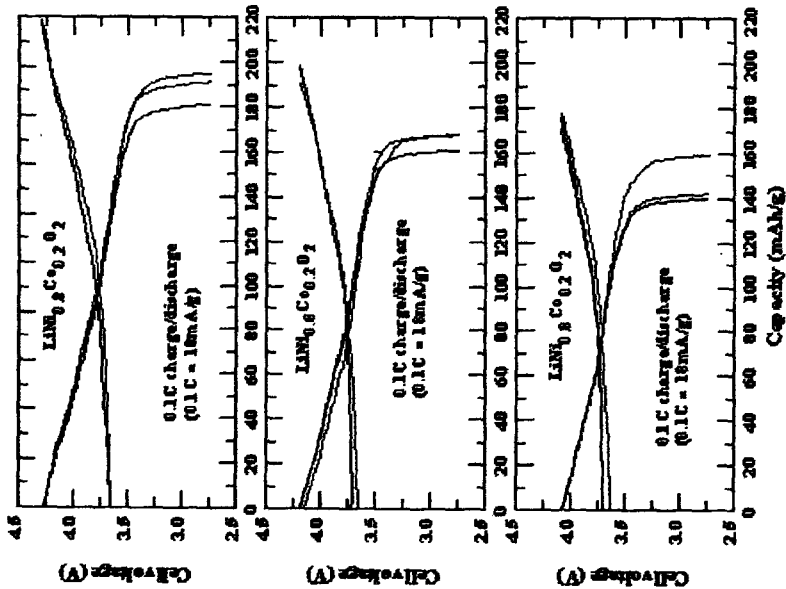


**Method B**

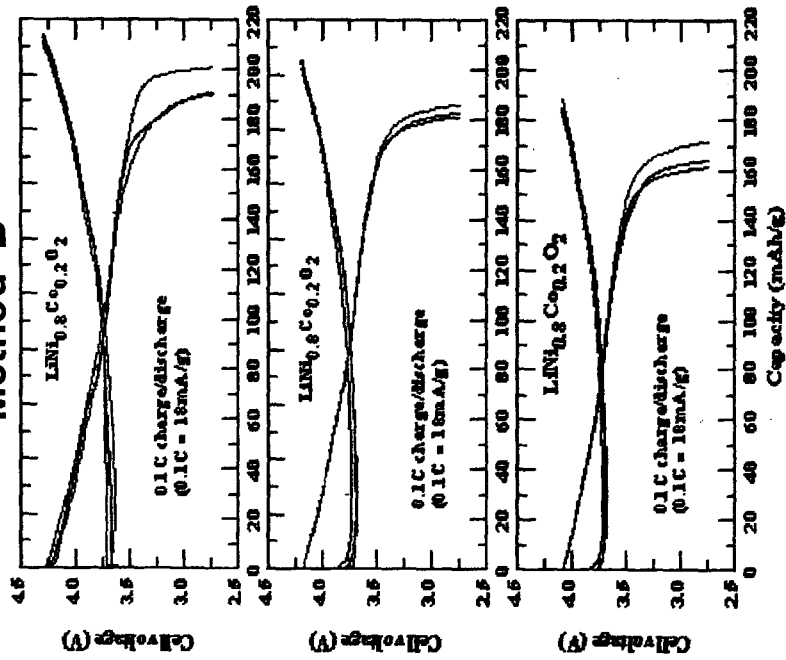


Comparison of the Cycling Curves in the First Cycle at Different Cut-off Voltages

Method A

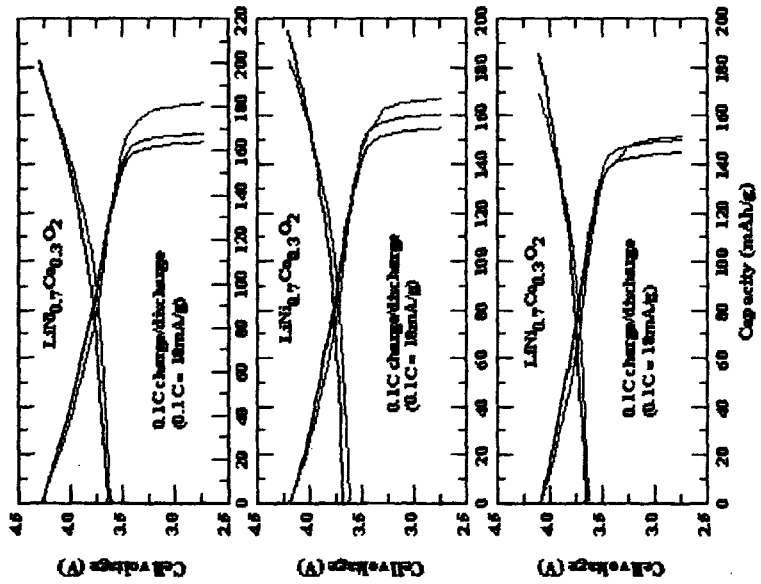


Method B

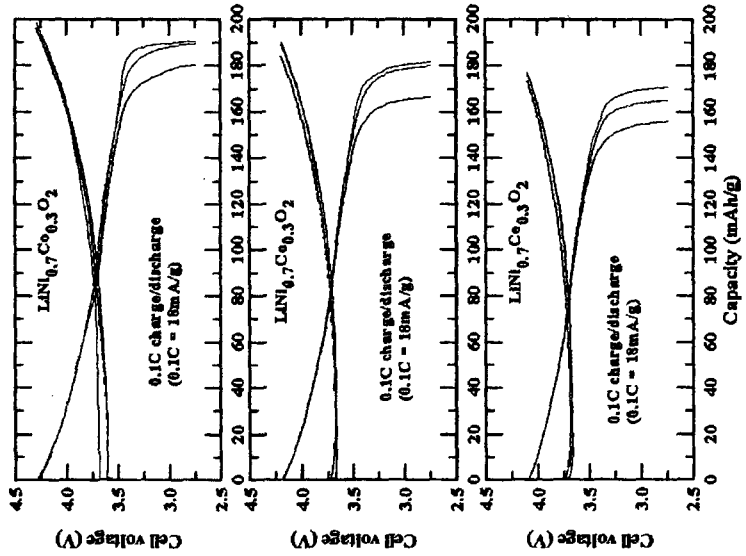


**Comparison of the Cycling Curves in the First Cycle at Different Cut-off Voltages**

**Method A**

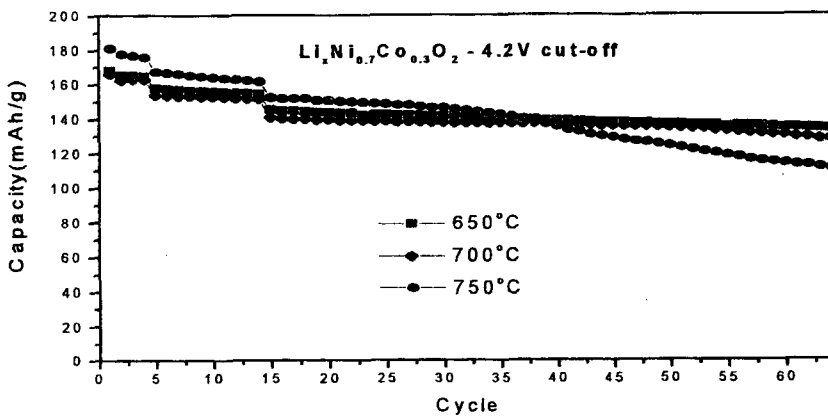
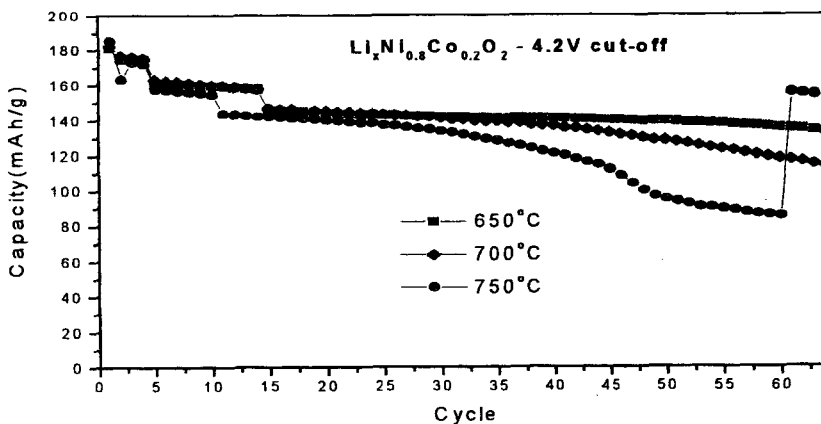
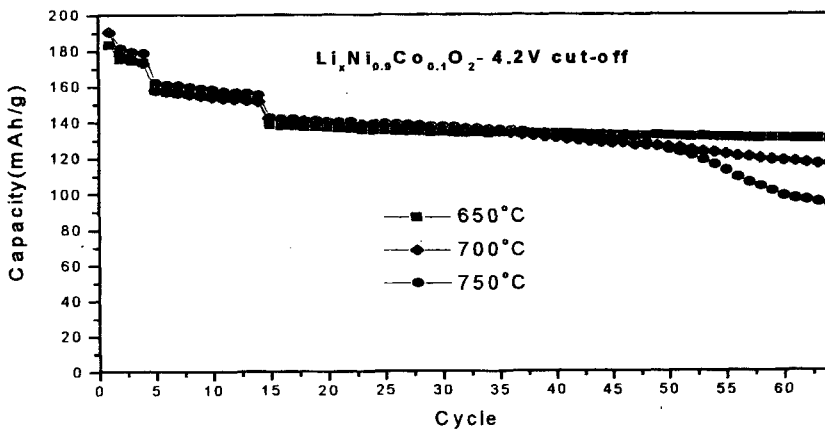


**Method B**





### Comparison of the Cyclability at Different Composition



## Conclusions

- $\text{LiNi}_{0.7}\text{Co}_{0.3}\text{O}_2$  Prepared by Method B Showed the Highest Values of a/c in the  $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$  with  $x=0.1, 0.2$  and  $0.3$
- $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$  with Less Cubic Characteristics (i.e., Higher Anisotropy) Showed Better Cyclability Than Those with Higher One  $\rightarrow$  Increased Layered Characteristics
- $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$  Prepared by Method B Showed higher Capacity Retention and Better Cyclability Than Those Prepared by Method A
- $\text{LiNi}_{0.7}\text{Co}_{0.3}\text{O}_2$  Showed the Best Cycling Performance at Different Cut-off Voltages in the  $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$  with  $x=0.1, 0.2$  and  $0.3$  Prepared by Method B