

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

삼성전관 기술본부

에너지LAB

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

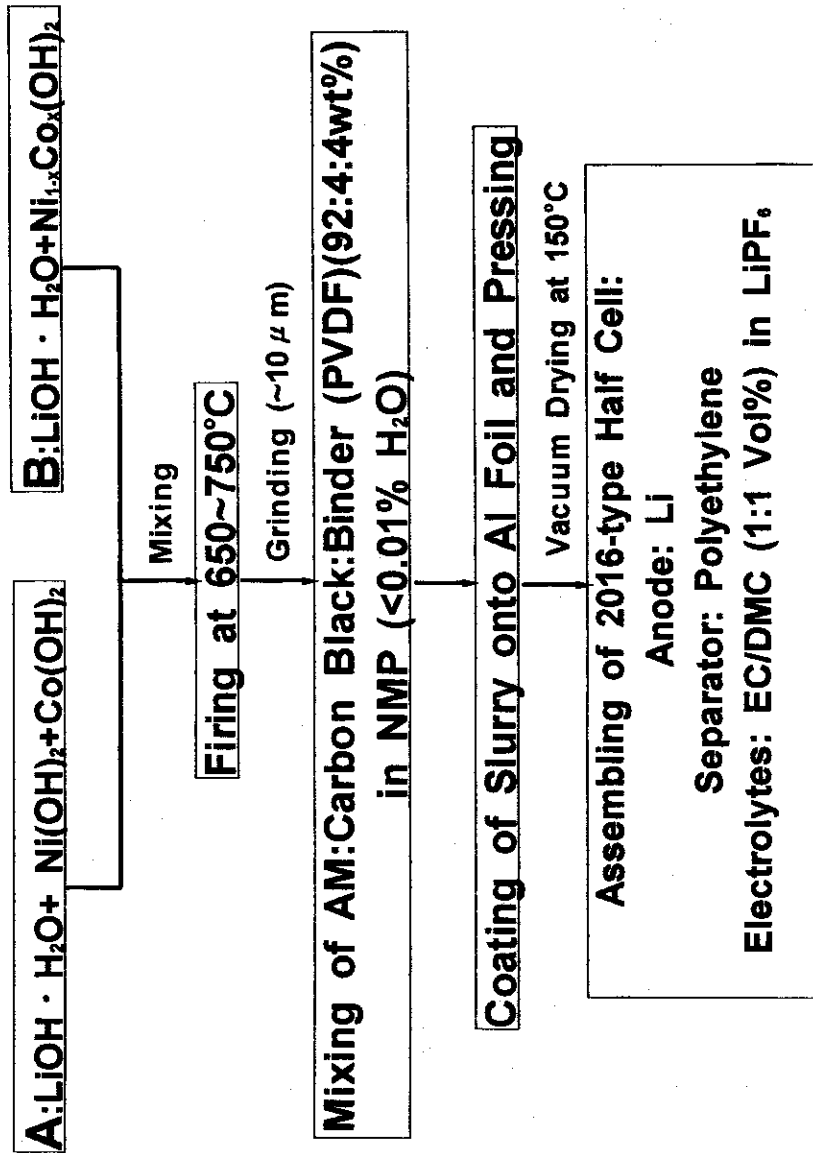
1. Optimization of Synthesis Method

- No Formation of 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 ($\approx 180\text{mA/g}$)
- Low Irreversible Capacity ($< 20\text{mAh/g}$)

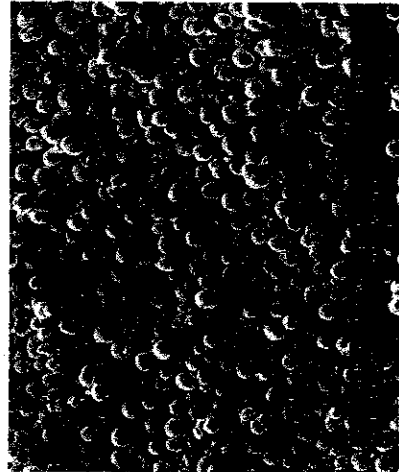
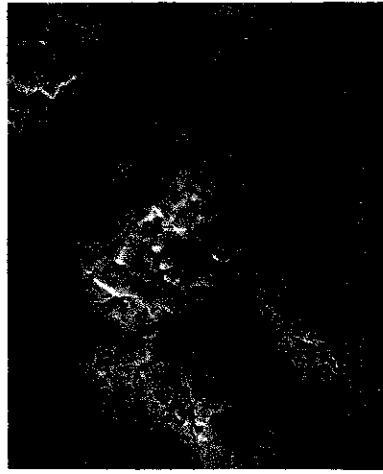
Experimental



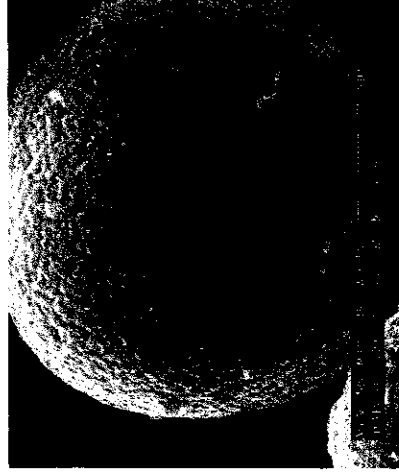
SEM Pictures of $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$ Prepared by Different Methods



Method A
Tap density:
2.6cc/g

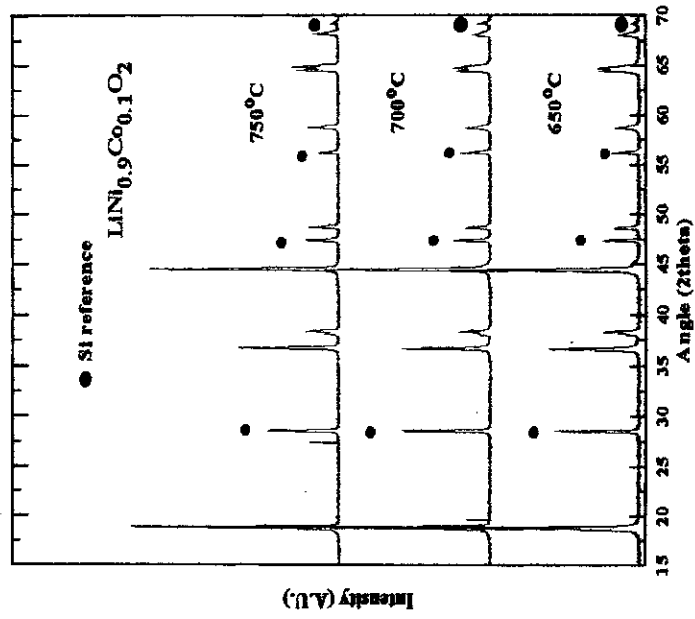


Method B
Tap density:
2.9cc/g

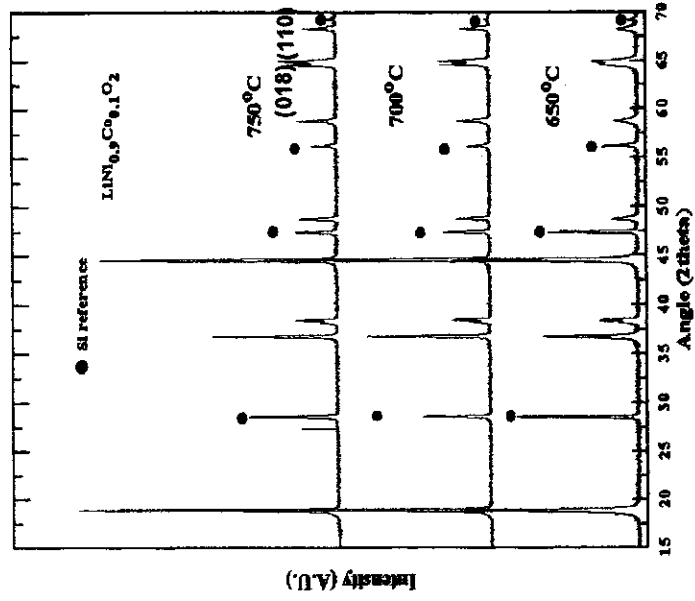


Comparison of the Powder XRD Patterns

Method A

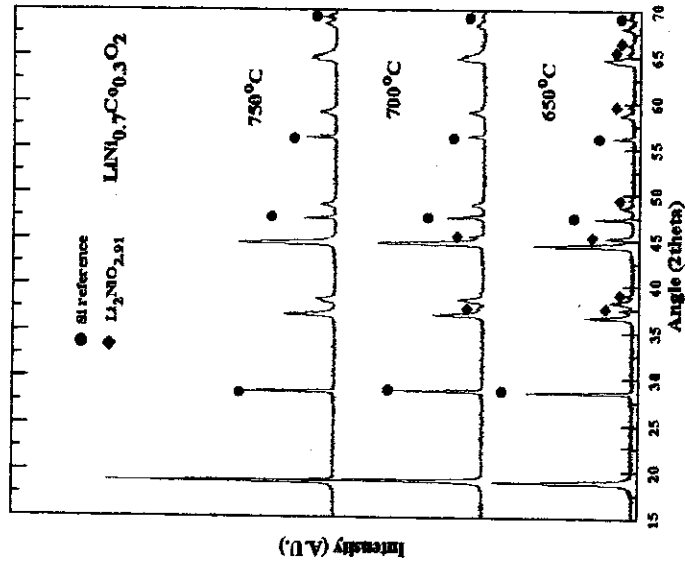


Method B

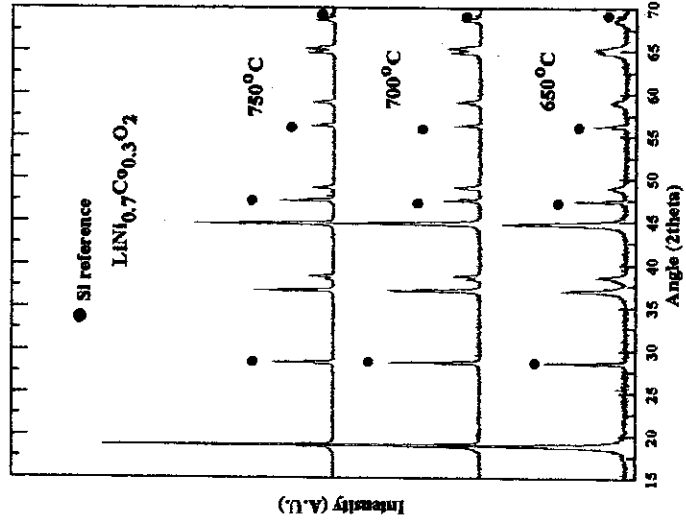


Comparison of the Powder XRD Patterns

Method A

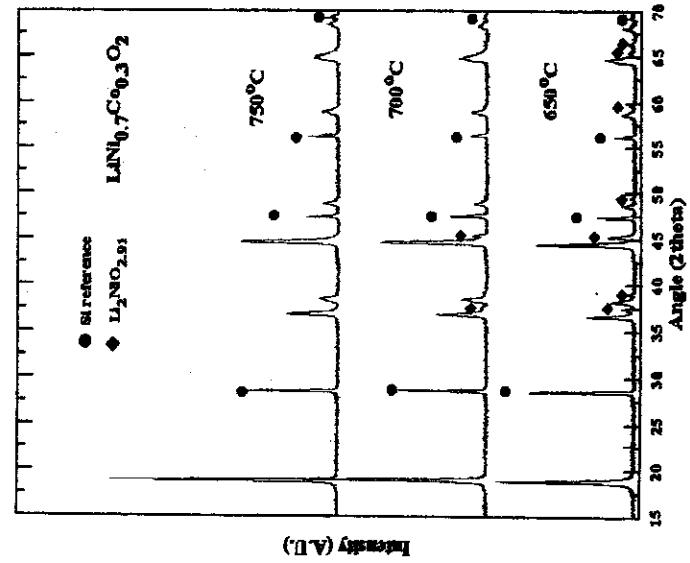


Method B

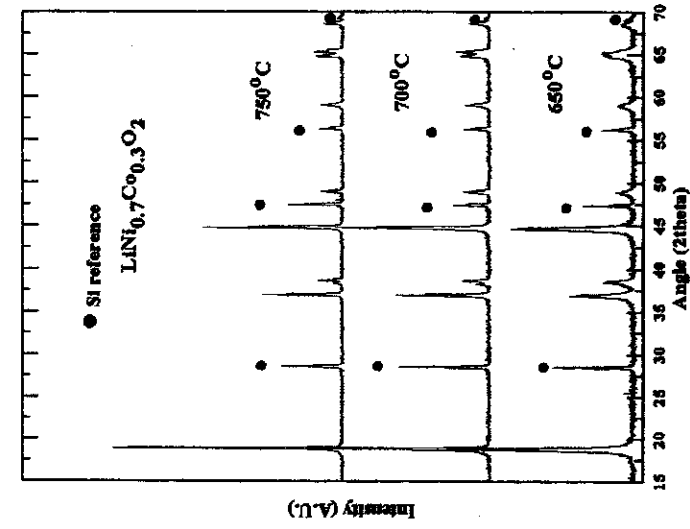


Comparison of the Powder XRD Patterns

Method A



Method B



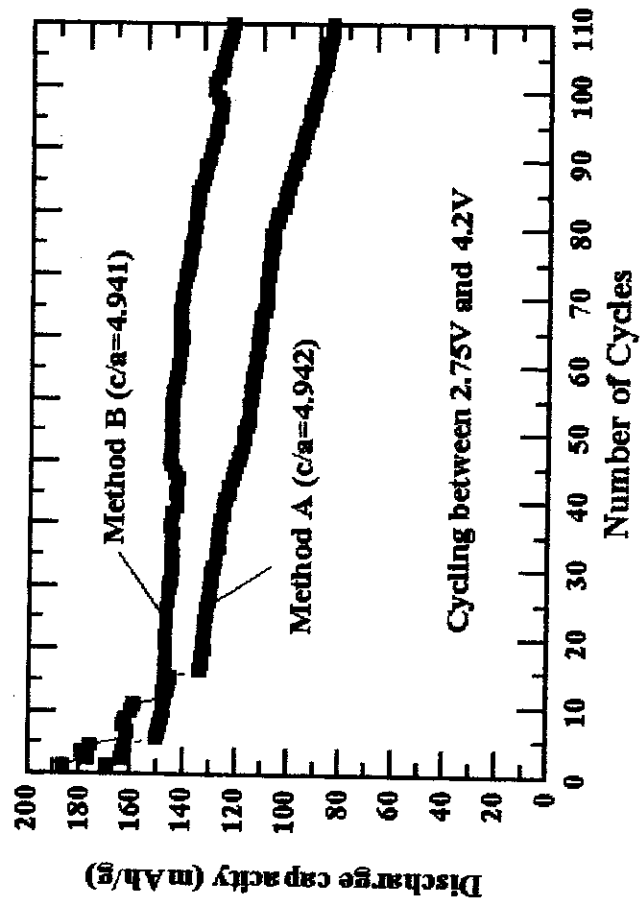
Comparison of a/c Values in the $\text{LiNi}_{1-x}\text{Co}_x\text{O}_2$ Materials Depending on Firing Temperature

Method	A			B		
	$\text{LiNi}_0.9\text{Co}_{0.1}\text{O}_2$	$\text{LiNi}_{0.8}\text{Co}_{0.2}\text{O}_2$	$\text{LiNi}_{0.7}\text{Co}_{0.3}\text{O}_2$	$\text{LiNi}_{0.6}\text{Co}_{0.4}\text{O}_2$	$\text{LiNi}_{0.5}\text{Co}_{0.5}\text{O}_2$	$\text{LiNi}_{0.4}\text{Co}_{0.6}\text{O}_2$
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 [#]	4.948

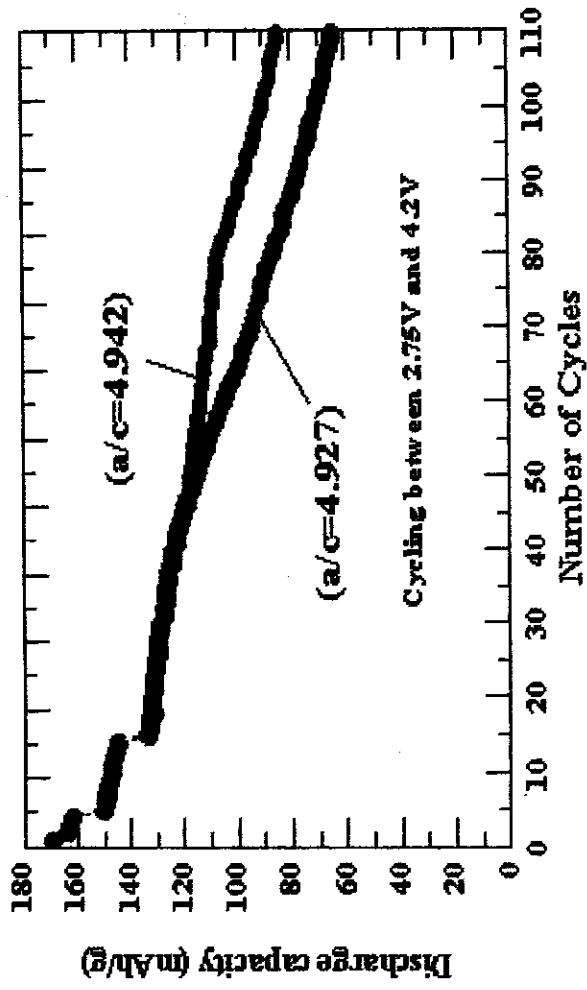
* LiCoO_2 : 4.99, LiNiO_2 : 4.935

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

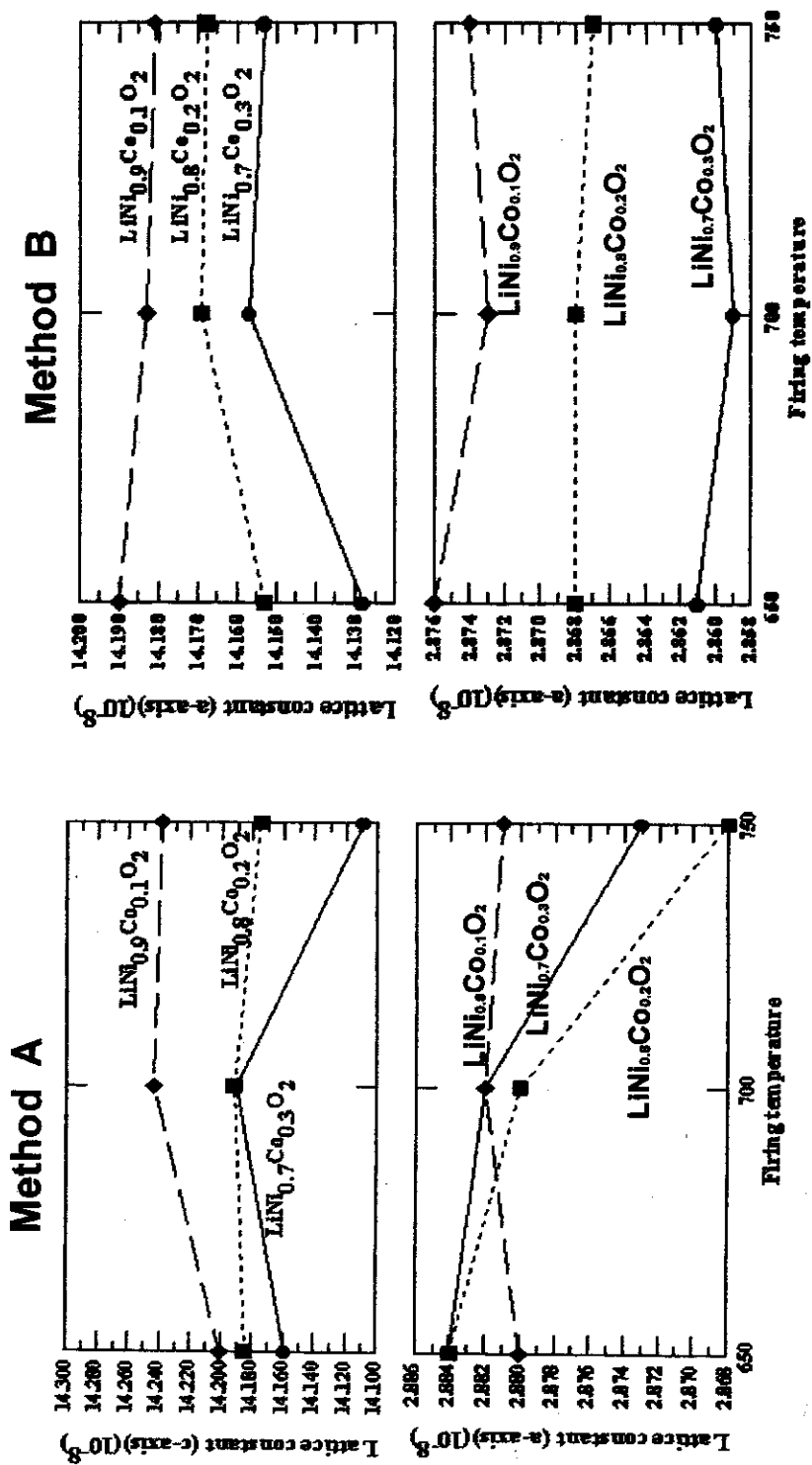
Cycling Behavior of $\text{LiNi}_{0.8}\text{Co}_{0.2}\text{O}_2$ Prepared by Different Methods



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

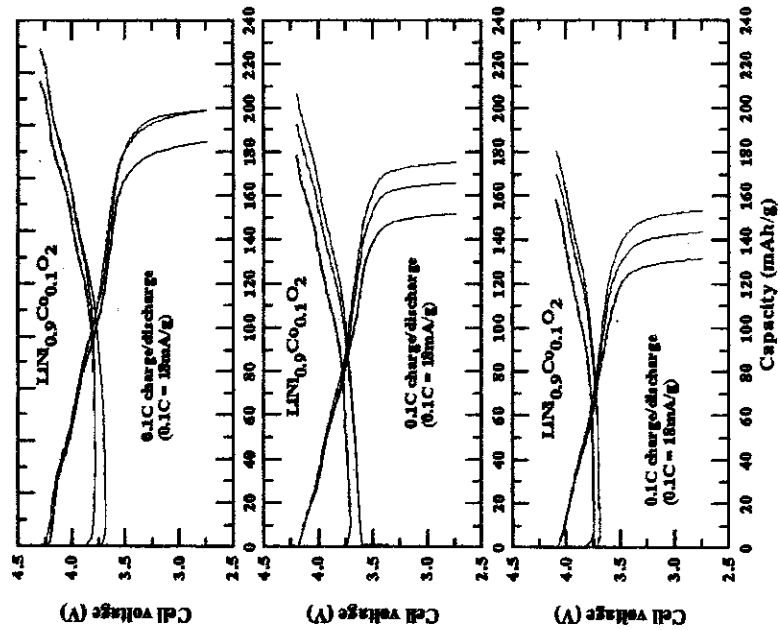


Comparison of Lattice Parameters as a Function of Firing Temperature

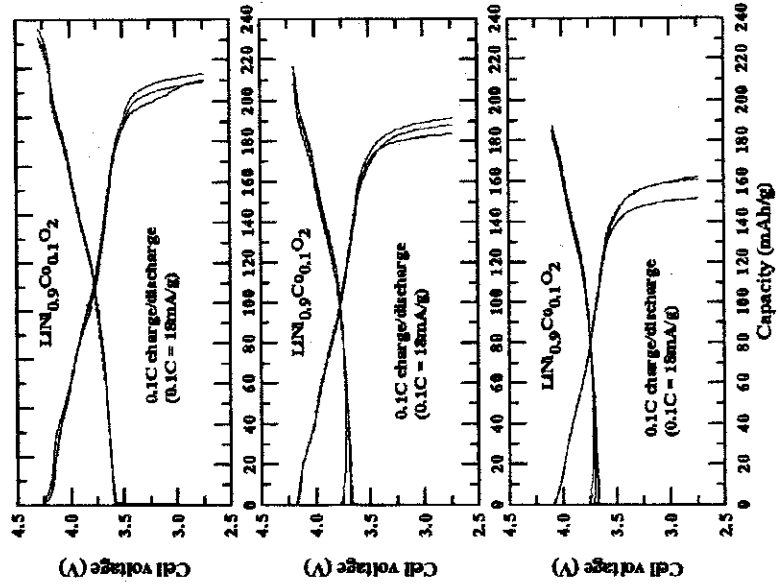


Comparison of the Cycling Curves in the First Cycle at Different Cut-off Voltages

Method A

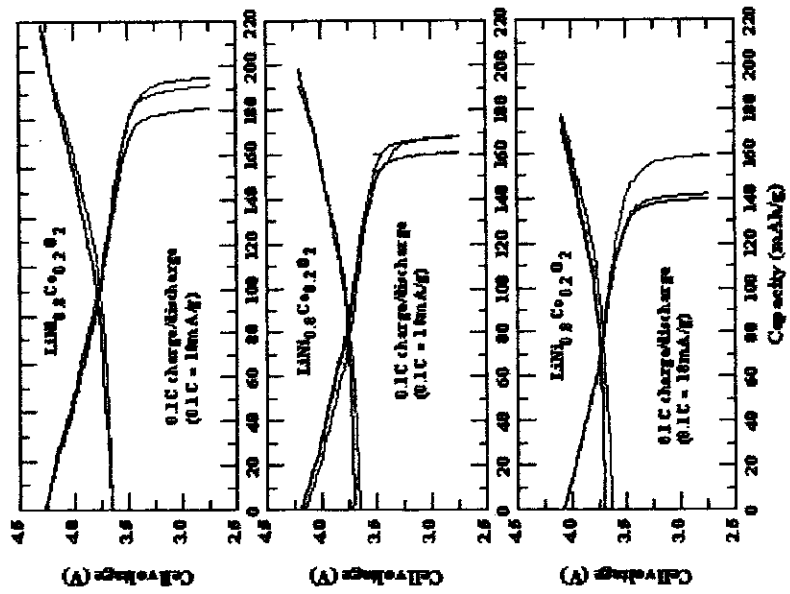


Method B

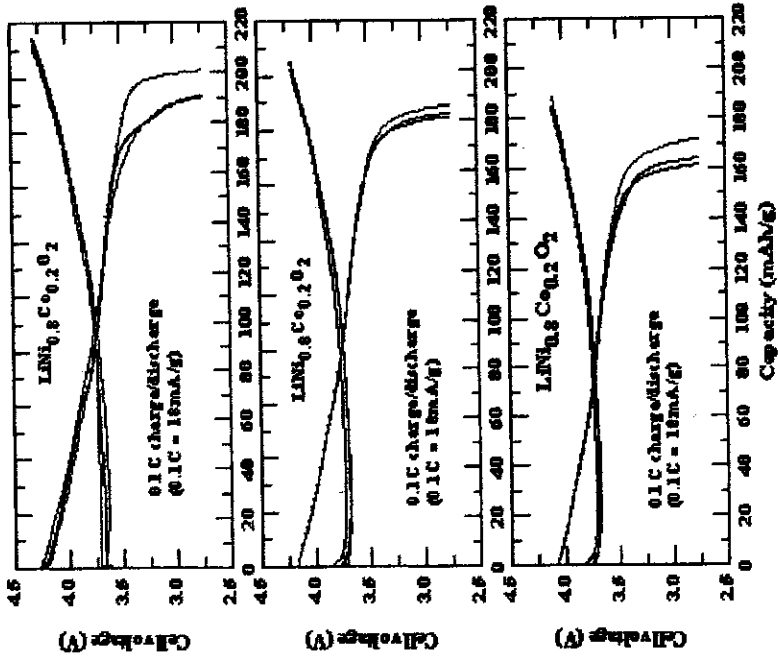


Comparison of the Cycling Curves in the First Cycle at Different Cut-off Voltages

Method A

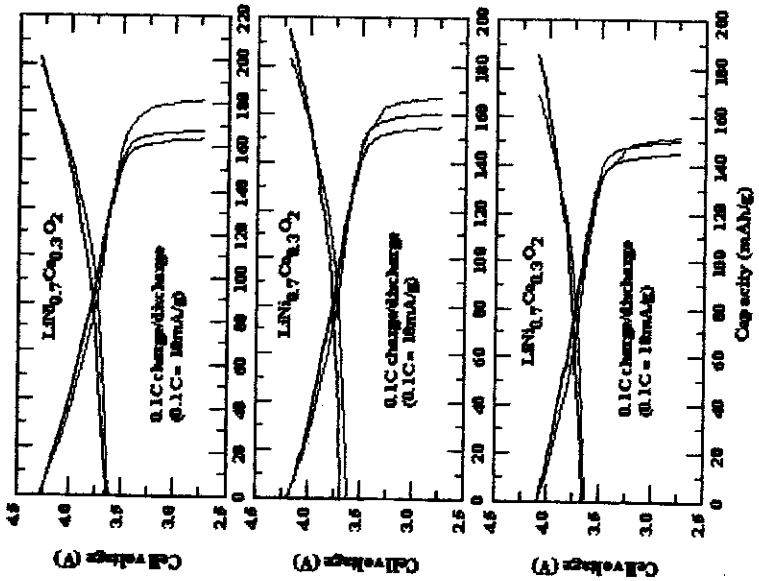


Method B

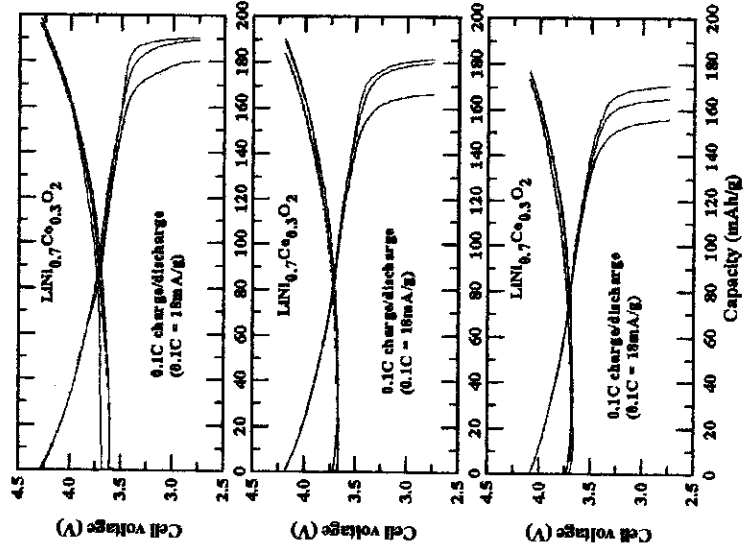


Comparison of the Cycling Curves in the First Cycle at Different Cut-off Voltages

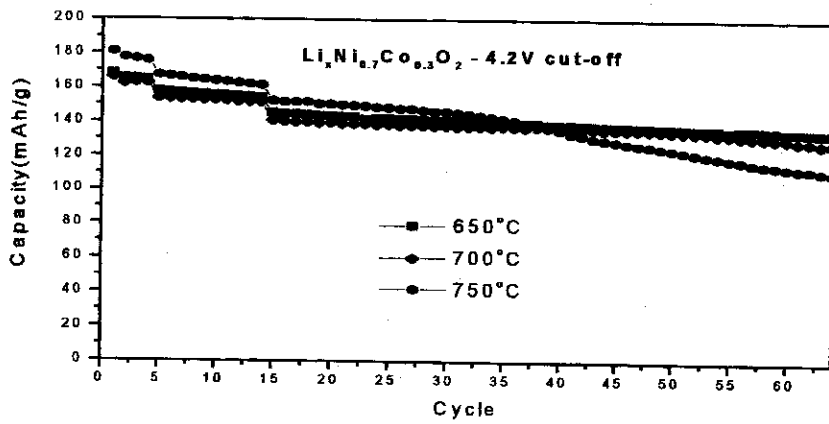
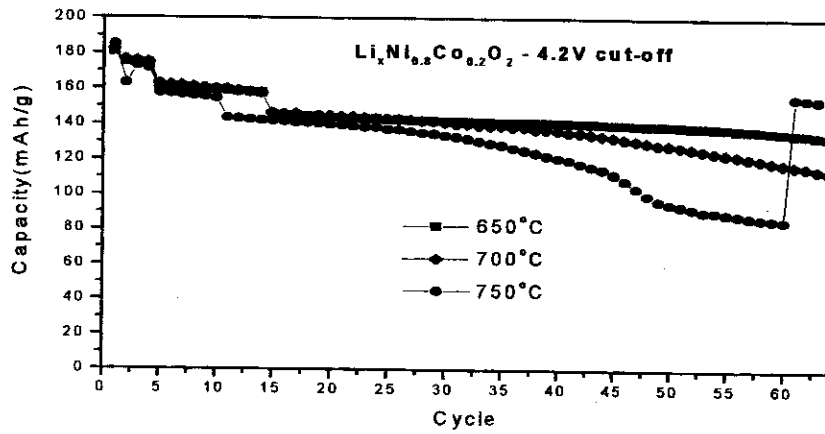
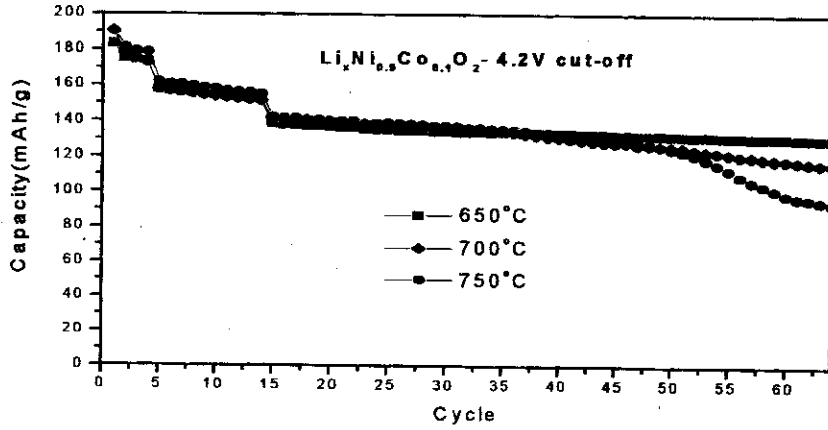
Method A



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