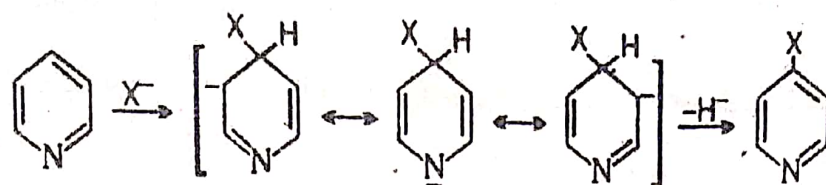


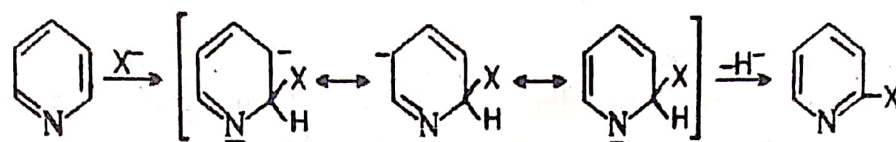
the substitution at position-2 and 4 in pyridine gives the similar resonating structures. After observing minutely the role of N-atom as -I effect, it attracts electron from position-2 more than that from 4.

**Nucleophilic attack on position-4 :**



Most Stable

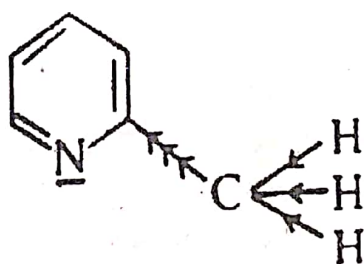
**Nucleophilic attack on position-2 :**



Most Stable

**Q.26. Methyl group in 2-position is very reactive in pyridine—Why ?**

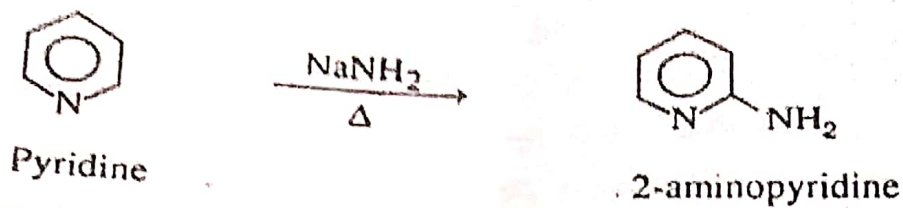
**Ans. :** Due to electronegative nature, N-atom becomes negative as a result of +M effect and C-atom of  $-CH_3$  group at 2-position becomes positive. Further  $-CH_3$  group has also +I effect which helps the +M effect to operate as shown in figure. This causes a positive charge on C-atom and thus C-H bonds are weakened and H-atom becomes reactive due to its labile nature—



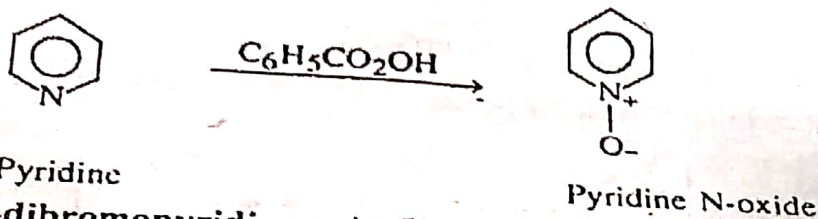
**Q.27. How would you convert pyridine into**

- (i) 2-amino pyridine
- (ii) Pyridine N-oxide
- (iii) 3, 5-dibromo pyridine.

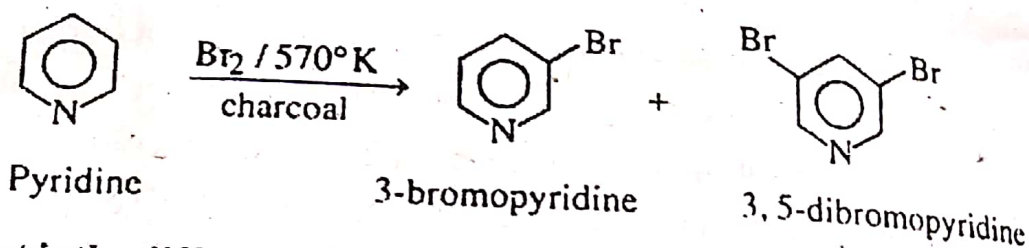
**Ans. :** (i) 2-amino pyridine : When pyridine is heated with sodamide, the main product 2-amino pyridine is formed—



(ii) **Pyridine N-oxide** : Pyridine is oxidised by  $\text{H}_2\text{O}_2$  or per acid to its N-oxide—

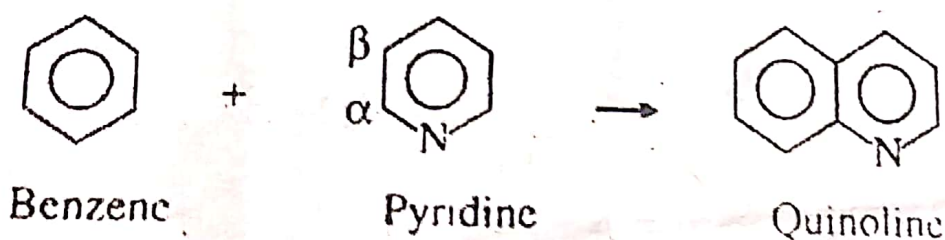


(iii) **3, 5-dibromopyridine** : At 570K in the presence of charcoal or pumice, pyridine forms a mixture of 3-bromopyridine and 3, 5-dibromopyridine—



**Q.28. What is the difference between quinoline and isoquinoline?**

**Ans. :** Quinoline has a benzene ring fused to  $\alpha, \beta$ - positions of a pyridine ring. So, it is  $\alpha, \beta$  benzopyridine—



whereas isoquinoline has a benzene ring fused to  $\beta, \gamma$  - positions of pyridine ring. So it is  $\beta, \gamma$  benzopyridine—

