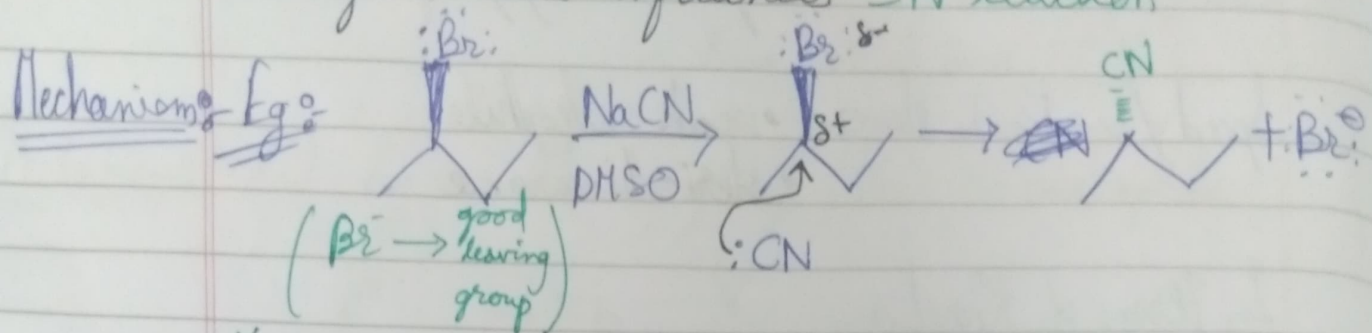


Reaction: In S_N2 reactions, a catalyst-like H^+ is not required. This is because, the nucleophile attacks the molecule directly.

There is no need for a catalyst as the nucleophile's strength is what influences S_N2 reaction.



Here, $\text{NaCN} \rightarrow \text{Na}^+ + \text{CN}^-$
strong nucleophile

DMSO \rightarrow Polar Aprotic solvent

From above reaction,

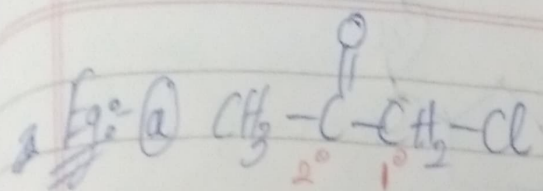
The nucleophile CN^- attacks the molecule on rear side

The Backside of the molecule is attacked, it causes an inversion of the molecule after the leaving group leaves & nucleophile gets attached.

This Inversion is called Walden Inversion

Steric Hindrance However, Transition State's Stability is also influenced by Steric Hindrance

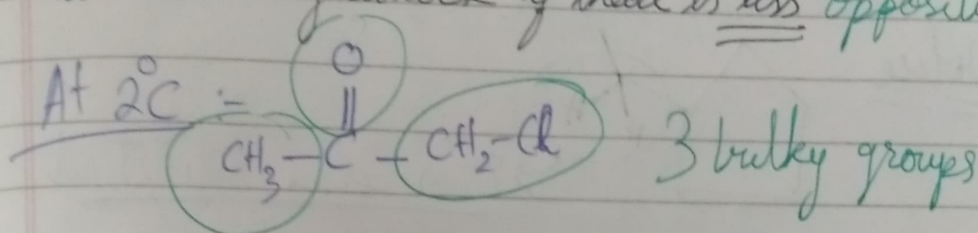
It is the amount of bulky groups surrounding C which causes difficulty for ~~an~~ nucleophile to attack.



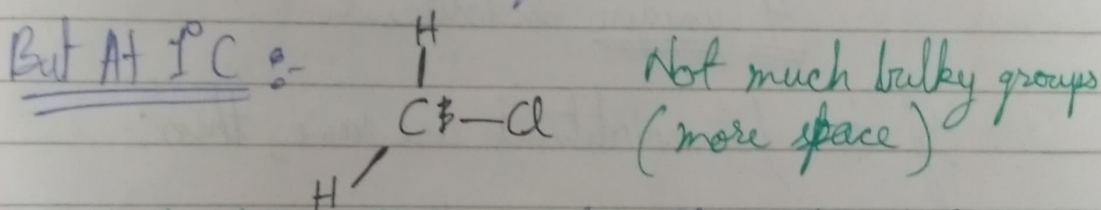
Here, the 1°C is the one with the Lg [leaving group]
 2°C is the next carbon

Now generally, $3^\circ > 2^\circ > 1^\circ$
Stability

But, for a nucleophile to attack molecule, it can only attack if there is less opposition



\therefore Nucleophile will be less favourable to attack at 2°C position



\therefore Nucleophile will be more favourable to attack at 1°C position

\therefore Steric hinderance is less when there is lesser bulkier groups attached to carbon.

\therefore For SN^2 : $1^\circ > 2^\circ > 3^\circ$ [Steric]
 \leftarrow Steric hinderance decreases

$\therefore 1^\circ \text{C}$ is most favoured by nucleophiles

In 2° & 3°C , steric hinderance increases

$$\text{Steric Hinderance} \propto \frac{1}{\text{SN}^2}$$