

# Summary

1. Substitution reactions at 1° and 2° (but not 3°) C(sp<sup>3</sup>) centers usually proceed by the S<sub>N</sub>2 mechanism under basic or neutral conditions. Substitution at C(sp<sup>2</sup>) centers rarely (if ever) occurs by the S<sub>N</sub>2 mechanism.
2. Leaving group ability:
  - The weaker the base, the better is the leaving ability.
  - F<sup>-</sup>, OH<sup>-</sup>, RO<sup>-</sup> (except epoxides), H<sup>-</sup>, and carbanions are almost never leaving groups in [S<sub>N</sub>1] or [S<sub>N</sub>2] reactions.
3. Nucleophilicity (in an **aprotic** solvent like DMSO):
  - The stronger the base, the stronger the nucleophile.
4. Nucleophilicity (in a **protic** solvent like methanol):
  - Neutrals are poorer nucleophiles than anions.
  - Comparing nucleophilic atoms that are in the same **row**, the stronger the base, the stronger the nucleophile.
  - Comparing nucleophilic atoms that are in the same **column**, the larger the atom, the stronger the nucleophile.