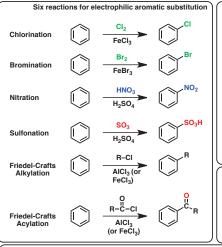
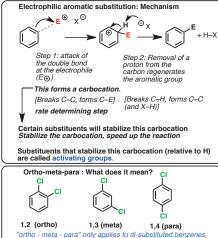
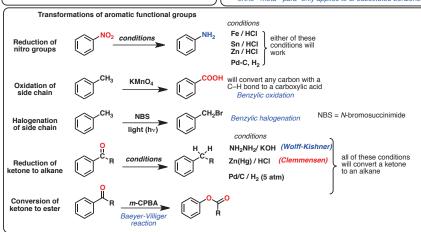
Reactions of Aromatic Compounds

"Master Organic Chemistry" masterorganicchemistry.com

Note - this sheet is not meant to be comprehensive. Your course may provide additional material, or may not cover some of the reactions shown here. Your course instructor is the final authority.







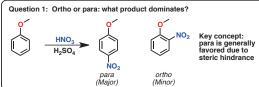
The group X on the ring will affect the stability of the carbocation X B Ortho The group X on the ring will affect the stability of the carbocation X H Ortho E Ortho F Ortho

What if there's already a group on the ring?

Activating vs. Deactivating: What does it mean?

aromatic substitution to be faster, relative to H A deactivating group is a group that causes electrophilic aromatic substitution to be slower, relative to H CH₃ CI CO₂Me relative rate for nitration: CH₃ is activating

An activating group is a group that causes electrophilic



Resonance forms A and B are KEY

If **X** has an electron pair, these resonance forms will be **stabilized** through formation of a double bond:

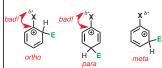
This resonance form is available for the ortho and para adducts, but NOT the meta.

This is why the following groups are ortho-para "directors"

Alkyl groups such as CH₃, CH₂CH₃, etc. are also ortho-para directors. Why? Notice the partial charge

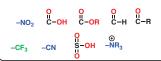
 H^{δ^*} The partial negative charge on the carbon $-L^{\delta^*}$ will stabilize an adjacent positive charge. $-L^{\delta^*}$

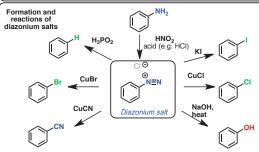
Similarly, if there is a partial positive charge on the atom adjacent to the ring, this will **destabilize** resonance forms A and B

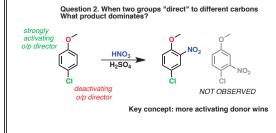


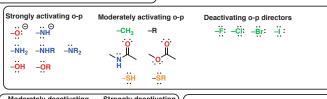
Destabilized! adjacent Not as unstable positive charges

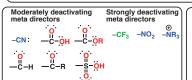
This is why the following groups are meta directors. (Although "ortho-para avoider" is more appropriate)











CI and CO₂Me are deactivating

Omissions, Mistakes, Suggestions?

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