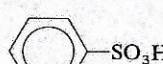
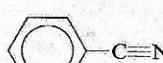
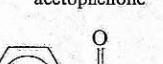
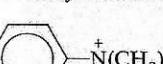


~order of deactivating "power":

halogens  $< \text{---C---OR}$   $< \text{---C---OH}$   $< \text{---C---R}$   $< \text{---C---H}$   $< \text{SO}_3\text{H}$   $< \text{C}\equiv\text{N}$   $< \text{NO}_2$   $< \text{NR}_3^+$

### DEACTIVATING META-DIRECTORS

Group	Resonance Structures	Example
$-\text{NO}_2$ nitro	$\left[ -\text{N}^+ \text{---} \begin{matrix} \ddot{\text{O}} \\   \\ \text{O} \end{matrix} \longleftrightarrow -\text{N}^+ \text{---} \begin{matrix} \ddot{\text{O}}^- \\   \\ \text{O} \end{matrix} \right]$	 nitrobenzene
$-\text{SO}_3\text{H}$ sulfonic acid	$\left[ \begin{matrix} \ddot{\text{O}} \\   \\ \text{S} \\    \\ \text{O} \end{matrix} \text{---} \ddot{\text{O}}-\text{H} \longleftrightarrow \begin{matrix} \ddot{\text{O}} \\   \\ \text{S} \\    \\ \text{O} \end{matrix} \text{---} \ddot{\text{O}}^-\text{H} \longleftrightarrow \begin{matrix} \ddot{\text{O}} \\   \\ \text{S} \\    \\ \text{O} \end{matrix} \text{---} \ddot{\text{O}}^-\text{H} \right]$	 benzenesulfonic acid
$-\text{C}\equiv\text{N}$ : cyano	$\left[ -\text{C}\equiv\text{N} : \longleftrightarrow -\text{C}^+ \equiv \ddot{\text{N}} : \right]$	 benzonitrile
$-\text{C}(=\text{O})-\text{R}$ ketone or aldehyde	$\left[ \begin{matrix} \ddot{\text{O}} \\   \\ \text{C} \\    \\ \text{R} \end{matrix} \longleftrightarrow -\text{C}^+ \text{---} \ddot{\text{O}}^- \right]$	 acetophenone
$-\text{C}(=\text{O})-\text{O}-\text{R}$ ester	$\left[ \begin{matrix} \ddot{\text{O}} \\   \\ \text{C} \\    \\ \text{O} \end{matrix} \text{---} \ddot{\text{O}}-\text{R} \longleftrightarrow -\text{C}^+ \text{---} \ddot{\text{O}}-\text{R} \longleftrightarrow -\text{C}^+ \text{---} \ddot{\text{O}}^-\text{R} \right]$	 methyl benzoate
$-\text{NR}_3^+$ quaternary ammonium	$\begin{matrix} \text{---} \begin{matrix} \text{N}^+ \\   \\ \text{R} \\   \\ \text{R} \\   \\ \text{R} \end{matrix} \end{matrix}$	 trimethylanilinium iodide