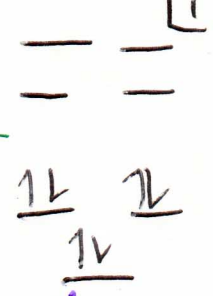
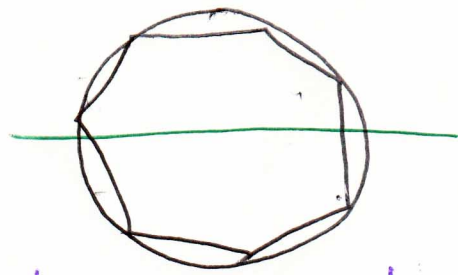
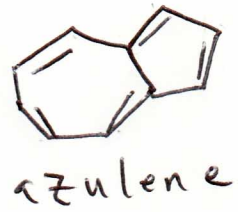
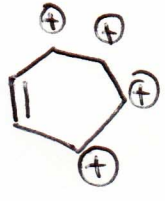


tropylium ion

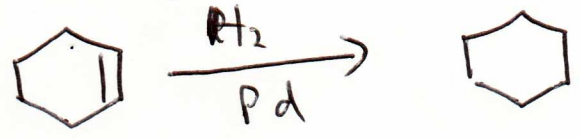


Bonding orbitals are exactly filled, so system is favorable

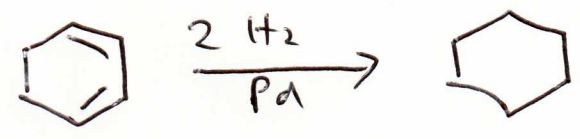


✓ aromatic (cyclic, planar, conjugated, 10 π e⁻)
For a Frost circle, a decagon would be inscribed.

Stability of benzene



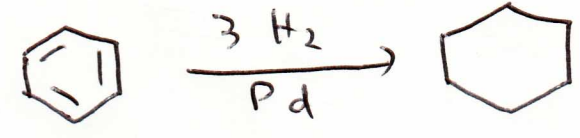
$\Delta H_{rxn} = -120 \text{ kJ/mol}$



$\Delta H_{rxn} (\text{predicted}) = -240 \text{ kJ/mol}$

$\Delta H_{rxn} (\text{actual}) = -232 \text{ kJ/mol}$

$|\Delta \Delta H_{rxn}| = 8 \text{ kJ/mol}$

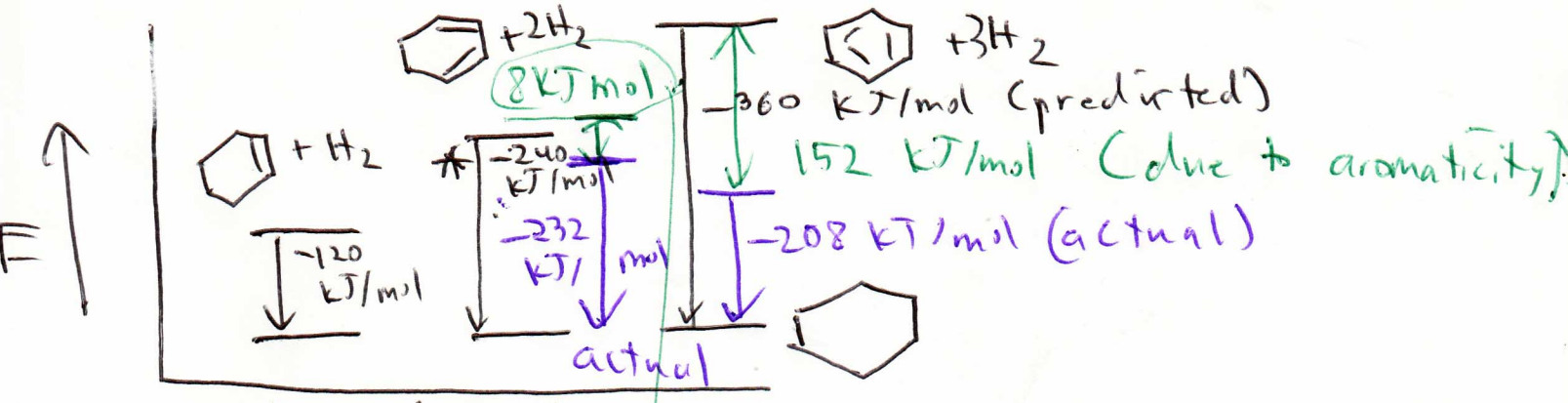


$\Delta H_{rxn} (\text{predicted}) = -360 \text{ kJ/mol}$

$\Delta H_{rxn} (\text{actual}) = -208 \text{ kJ/mol}$

$|\Delta \Delta H_{rxn}| = 152 \text{ kJ/mol}$

Quiz #3 3/7/12



* predicted

due to conjugation

The fact that benzene releases less energy than predicted means it is lower in energy to begin with ("stable").

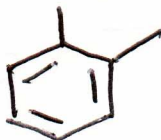
No nomenclature

Common names

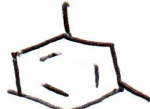
(ortho, meta, para)



toluene



o-xylene



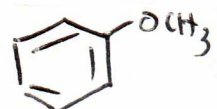
m-xylene



p-xylene



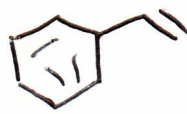
phenol



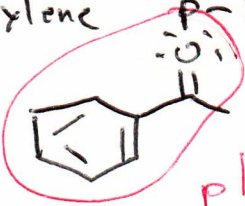
anisole



benzaldehyde



styrene



phenone

acetophenone



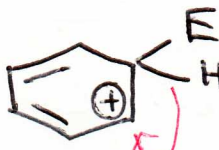
1-ethyl-3-methylbenzene

Whenever a molecule can be numbered equally in two ways, the lowest # is given to the substituent that is lower alphabetically

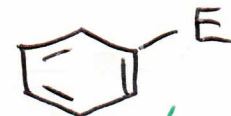
Electrophilic Aromatic Substitution



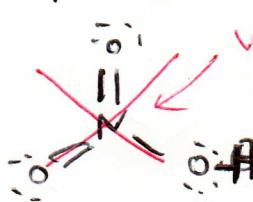
E^+



E
 H
base



Nitration



violates octet rule

