# Chapter 24 Phenols

## 24.1 Nomenclature

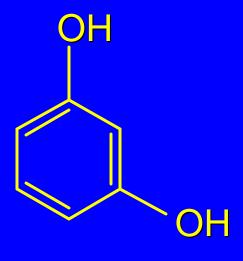
#### Nomenclature



named on basis of phenol as parent substituents listed in alphabetical order

lowest numerical sequence: first point of difference rule

#### Nomenclature





1,2-Benzenediol

(common name: pyrocatechol)

1,3-Benzenediol

(common name: resorcinol)

1,4-Benzenediol

(common name: hydroquinone)

#### Nomenclature



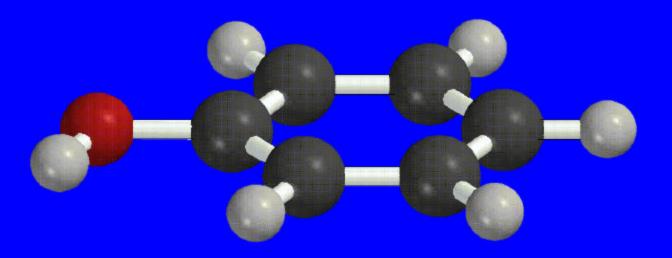
p-Hydroxybenzoic acid

name on basis of benzoic acid as parent

higher oxidation states of carbon outrank hydroxyl group

## 24.2 Structure and Bonding

#### Structure of Phenol



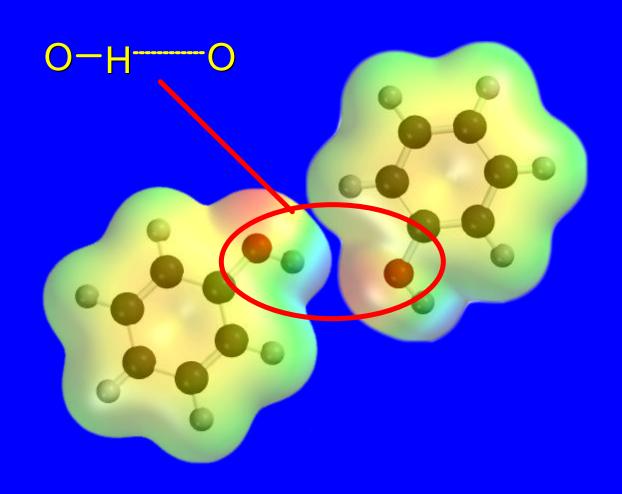
phenol is planar

C—O bond distance is 136 pm, which is slightly shorter than that of CH<sub>3</sub>OH (142 pm)

## 24.3 Physical Properties

The OH group of phenols allows hydrogen bonding to other phenol molecules and to water.

## Hydrogen Bonding in Phenols



### Physical Properties (Table 24.1)

Compared to compounds of similar size and molecular weight, hydrogen bonding in phenol raises its melting point, boiling point, and solubility in water.

## Physical Properties (Table 24.1)

	C <sub>6</sub> H <sub>5</sub> CH <sub>3</sub>	C <sub>6</sub> H <sub>5</sub> OH	C <sub>6</sub> H <sub>5</sub> F
Molecular weight	92	94	96
Melting point (°C)	<b>–95</b>	43	-41
Boiling point (°C,1 atm)	111	132	85
Solubility in H <sub>2</sub> O (g/100 mL,25°C)	0.05	8.2	0.2