

**TABLE 1.3** Less common yet important electrostatic interactions found in proteins.

Chemical species	Interaction	Interaction features
Aromatic rings	$\pi-\pi$	<ul style="list-style-type: none"> <li>• Involves two aromatic rings</li> <li>• Occurs between partially negative and partially positive regions in the rings</li> <li>• Highly geometry-dependent</li> </ul>
	cation- $\pi$	<ul style="list-style-type: none"> <li>• Often involve cationic metals</li> </ul>
	$\pi$ hydrogen bonds	<ul style="list-style-type: none"> <li>• The <math>\pi</math> electrons serve as the hydrogen bond acceptors</li> <li>• Weaker, yet geometrically more flexible than standard hydrogen bonds</li> </ul>
Sulfur and thiol	Hydrogen bonds	<ul style="list-style-type: none"> <li>• Involve the thiol (SH) group as a weak donor</li> <li>• May involve the two sulfur lone pairs as very weak acceptors</li> </ul>
	Thiol- $\pi$	<ul style="list-style-type: none"> <li>• Involve dipole-dipole interactions</li> <li>• May also involve interactions between the <math>\pi</math> electrons and the sulfur's low-lying <math>\sigma^*</math> orbitals</li> <li>• Highly geometry-dependent</li> </ul>
Lone electrons	$n \longrightarrow \pi^*$	<ul style="list-style-type: none"> <li>• Involves the lone electrons of electronegative atoms and the antibonding orbitals of other atoms</li> <li>• Estimated energy: <math>\sim 5\%</math>–<math>25\%</math> of a standard hydrogen bond</li> </ul>
Halogens	X-bonds	<ul style="list-style-type: none"> <li>• Electrostatic, occurring between a C–Cl/Br/I group and an electron-dense species (fully or completely charged)</li> <li>• Involve the internal dipole in the carbon-bound halogen</li> <li>• Strength depends on halogen: Cl &lt; Br &lt; I</li> </ul>