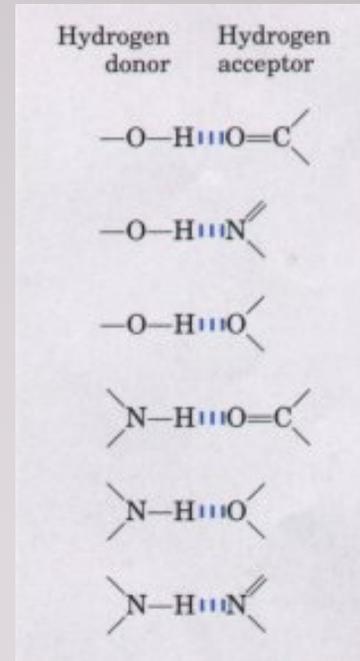
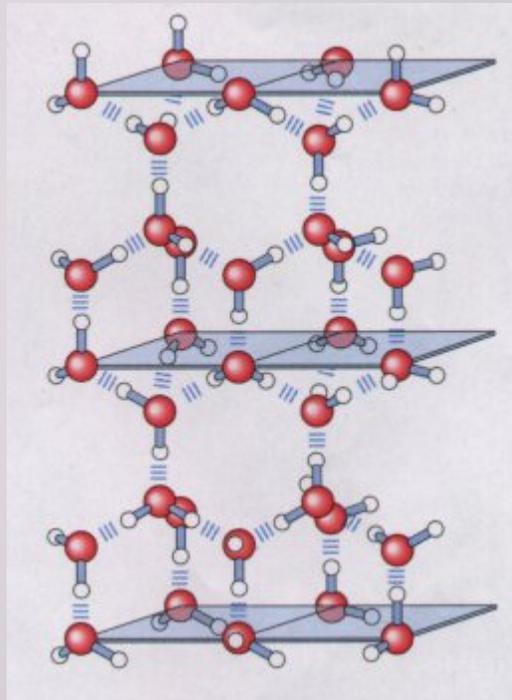
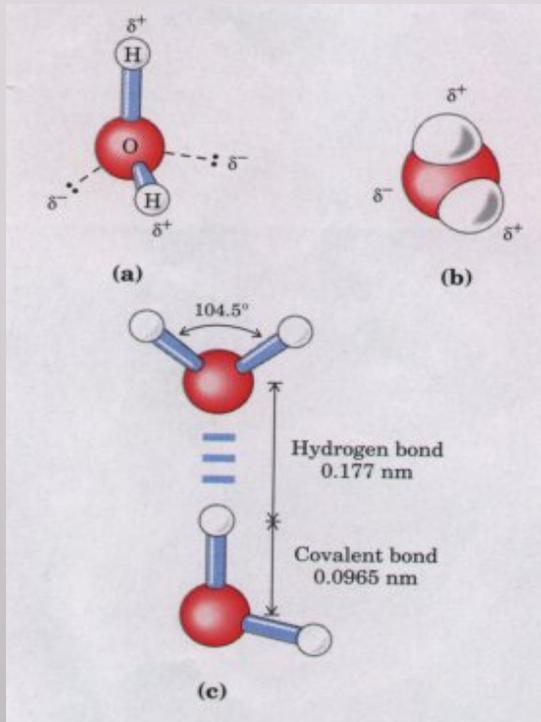


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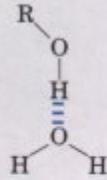
Table 4-1 Melting point, boiling point, and heat of vaporization of some common liquids

	Melting point (°C)	Boiling point (°C)	Heat of vaporization (J/g)*
Water	0	100	2,260
Methanol (CH ₃ OH)	-98	65	1,100
Ethanol (CH ₃ CH ₂ OH)	-117	78	854
Propanol (CH ₃ CH ₂ CH ₂ OH)	-127	97	687
Butanol (CH ₃ (CH ₂) ₂ CH ₂ OH)	-90	117	590
Acetone (CH ₃ COCH ₃)	-95	56	523
Hexane (CH ₃ (CH ₂) ₄ CH ₃)	-98	69	423
Benzene (C ₆ H ₆)	6	80	394
Butane (CH ₃ (CH ₂) ₂ CH ₃)	-135	-0.5	381
Chloroform (CHCl ₃)	-63	61	247

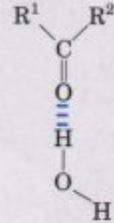
* The heat energy required to convert 1.0 g of a liquid at its boiling point, at atmospheric pressure, into its gaseous state at the same temperature. It is a direct measure of the energy required to overcome attractive forces between molecules in the liquid phase.



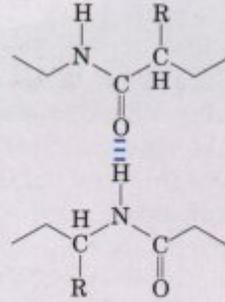
Between the hydroxyl group of an alcohol and water



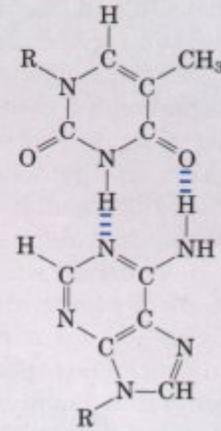
Between the carbonyl group of a ketone and water



Between two polypeptide chains

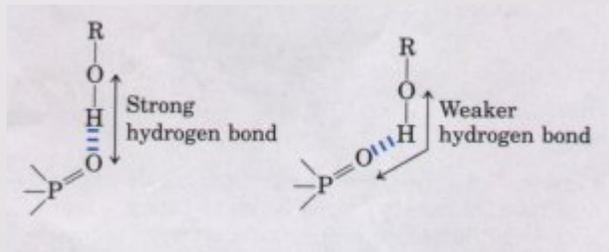


Between two complementary bases of two strands of DNA

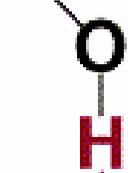


Thymine

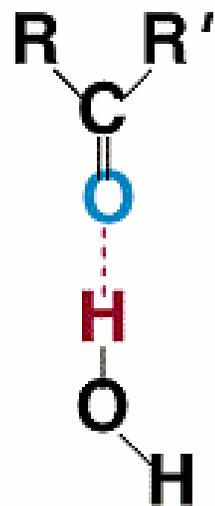
Adenine



Hydroxyl group

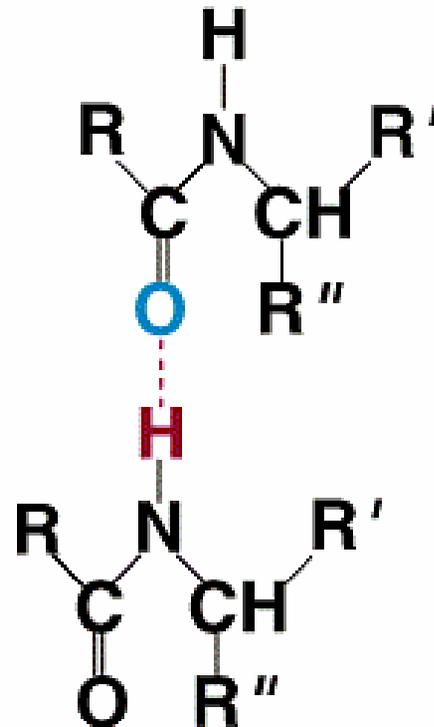


Water



Carbonyl

Water



Peptide chain

Peptide chain

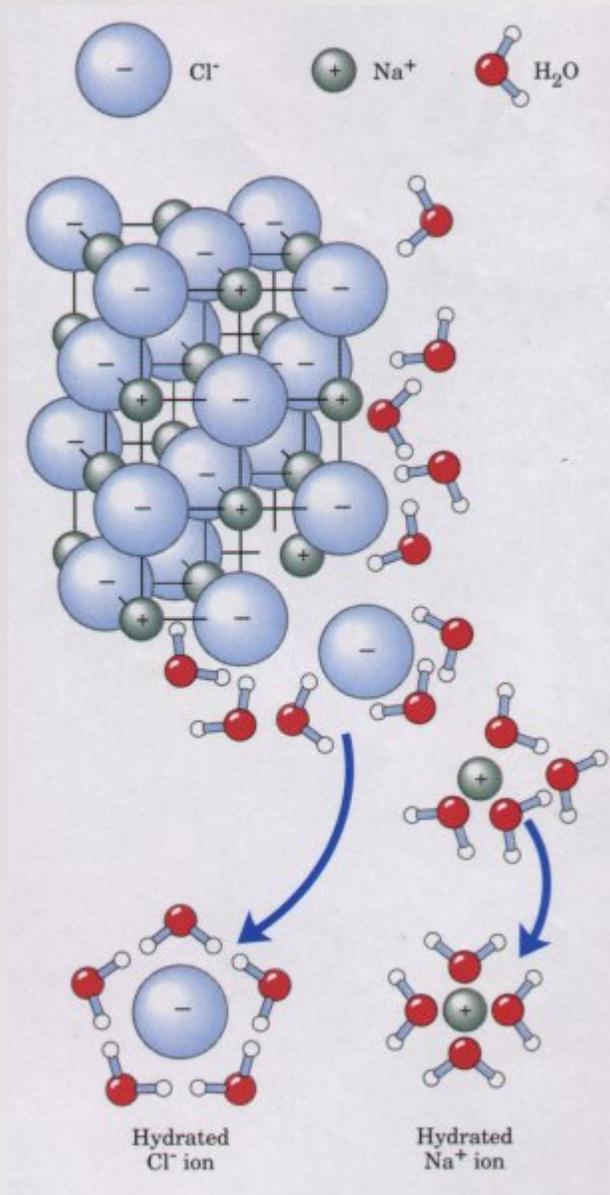


Table 4-3 Solubilities of some gases in water

Gas	Structure*	Polarity	Solubility in water (g/L)	Temperature (°C)
Nitrogen	$\text{N}\equiv\text{N}$	Nonpolar	0.018	40
Oxygen	$\text{O}=\text{O}$	Nonpolar	0.035	50
Carbon dioxide	$\overset{\delta-}{\text{O}}=\overset{\delta-}{\text{C}}=\overset{\delta-}{\text{O}}$	Nonpolar	0.97	45
Ammonia	$\begin{array}{c} \text{H} \\ \\ \text{H}-\text{N}-\text{H} \\ \\ \delta- \end{array}$	Polar	900	10
Hydrogen sulfide	$\begin{array}{c} \text{H} \\ \\ \text{H}-\text{S}-\text{H} \\ \\ \delta- \end{array}$	Polar	1,860	40

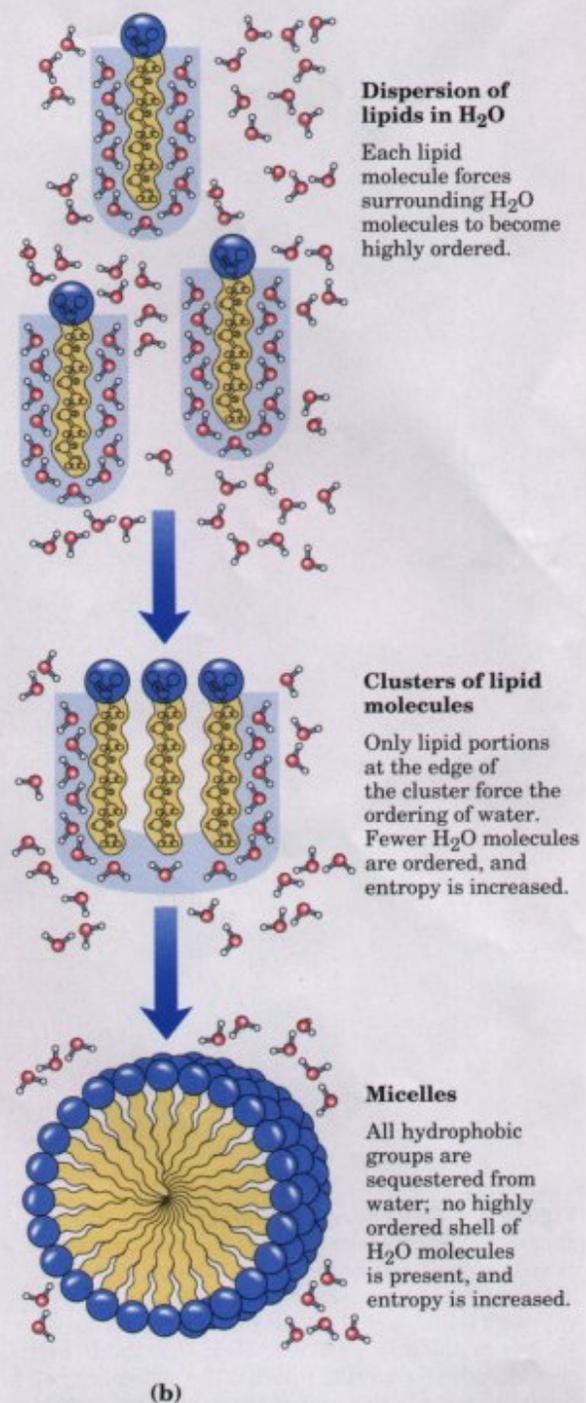
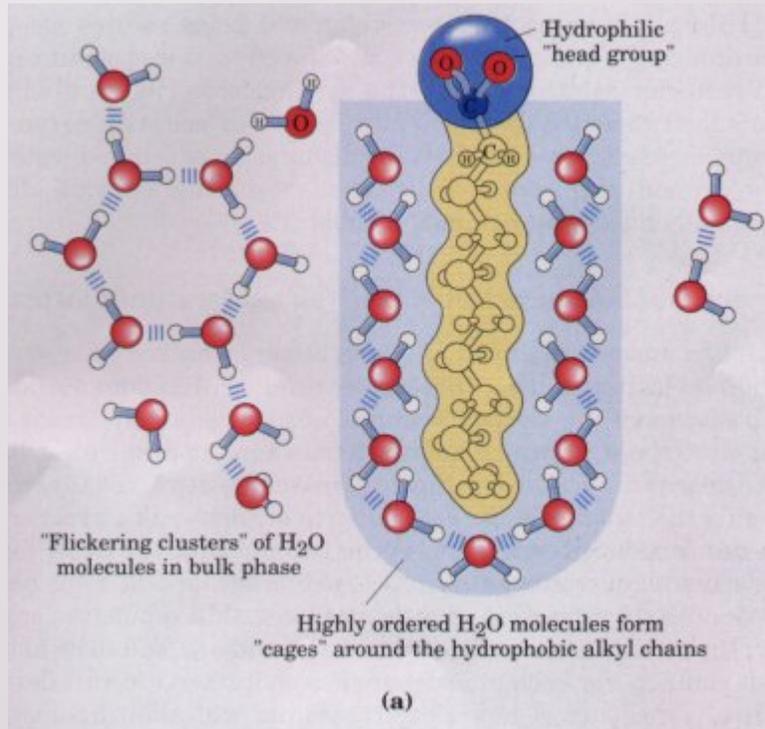


Table 4-4 Four weak interactions among biomolecules in aqueous solvent

Weak interaction		Stabilization energy (kJ/mol)
Hydrogen bonds		
Between neutral groups	$\diagup \text{C}=\text{O} \cdots \text{H}-\text{O}-$	8-21
Between peptide bonds	$\diagup \text{C}=\text{O} \cdots \text{H}-\text{N} \diagdown$	8-21
Ionic interactions		
Attraction	$-\text{NH}_3^+ \rightarrow \leftarrow \text{O}-\overset{\text{O}}{\parallel}{\text{C}}-$	42
Repulsion	$-\text{NH}_3^+ \leftrightarrow \text{H}_3\text{N}^+-$	≈ -21
Hydrophobic interactions	$\begin{array}{cc} \text{CH}_3 & \text{CH}_3 \\ & \diagdown \quad \diagup \\ & \text{CH} \\ & \\ & \text{CH}_2 \end{array} \quad \begin{array}{cc} \text{CH}_3 & \text{CH}_3 \\ & \diagdown \quad \diagup \\ & \text{CH} \\ & \\ & \text{CH}_2 \end{array}$	4-8
van der Waals interactions	Any two atoms in close proximity	4



$$K_{eq} = \frac{[\text{H}^+][\text{OH}^-]}{[\text{H}_2\text{O}]}$$

$$K_{eq} = \frac{[\text{H}^+][\text{OH}^-]}{55.5 \text{ M}}$$

$$55.5 \text{ M} \times K_{eq} = [\text{H}^+][\text{OH}^-] = K_w$$

$$K_{eq} \text{ agua} = 1.8 \times 10^{-16} \quad \text{A } 25^\circ\text{C}$$

$$55.5 \text{ M} \times 1.8 \times 10^{-16} = [\text{H}^+][\text{OH}^-]$$

$$1.0 \times 10^{-14} = [\text{H}^+][\text{OH}^-] = K_w$$

$$K_w = [\text{H}^+][\text{OH}^-] = [\text{H}^+]^2 \quad [\text{H}^+] = \sqrt{K_w} = \sqrt{1 \times 10^{-14}}$$

$$[\text{H}^+] = [\text{OH}^-] = 10^{-7} \text{ M}$$

$$\text{pH} = \log \frac{1}{[\text{H}^+]} = -\log [\text{H}^+]$$

Table 4-5 The pH scale

[H ⁺] (M)	pH	[OH ⁻] (M)	pOH*
10 ⁰ (1)	0	10 ⁻¹⁴	14
10 ⁻¹	1	10 ⁻¹³	13
10 ⁻²	2	10 ⁻¹²	12
10 ⁻³	3	10 ⁻¹¹	11
10 ⁻⁴	4	10 ⁻¹⁰	10
10 ⁻⁵	5	10 ⁻⁹	9
10 ⁻⁶	6	10 ⁻⁸	8
10 ⁻⁷	7	10 ⁻⁷	7
10 ⁻⁸	8	10 ⁻⁶	6
10 ⁻⁹	9	10 ⁻⁵	5
10 ⁻¹⁰	10	10 ⁻⁴	4
10 ⁻¹¹	11	10 ⁻³	3
10 ⁻¹²	12	10 ⁻²	2
10 ⁻¹³	13	10 ⁻¹	1
10 ⁻¹⁴	14	10 ⁰ (1)	0

* The expression pOH is sometimes used to describe the basicity, or OH⁻ concentration, of a solution; pOH is defined by the expression $\text{pOH} = -\log [\text{OH}^-]$, which is analogous to the expression for pH. Note that for all cases, $\text{pH} + \text{pOH} = 14$.

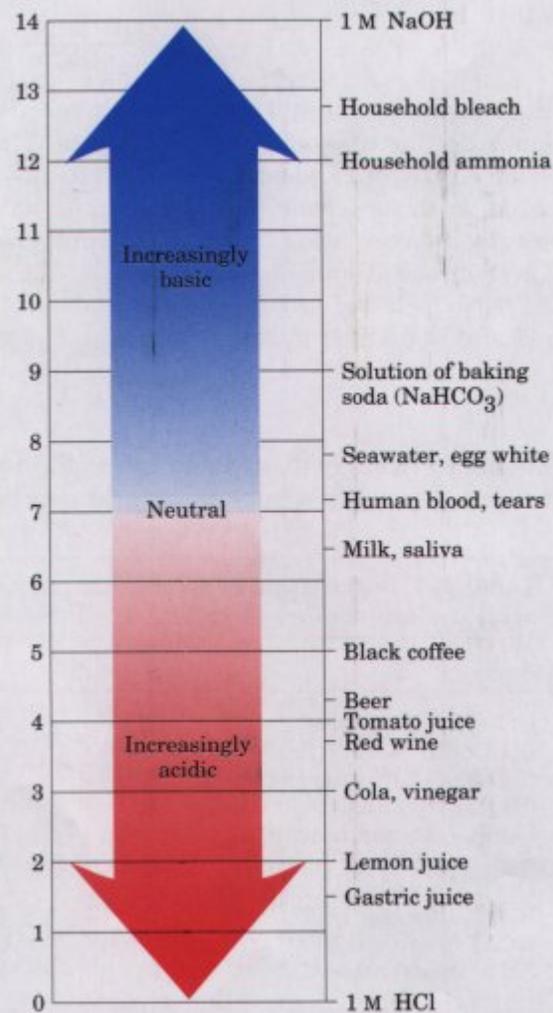
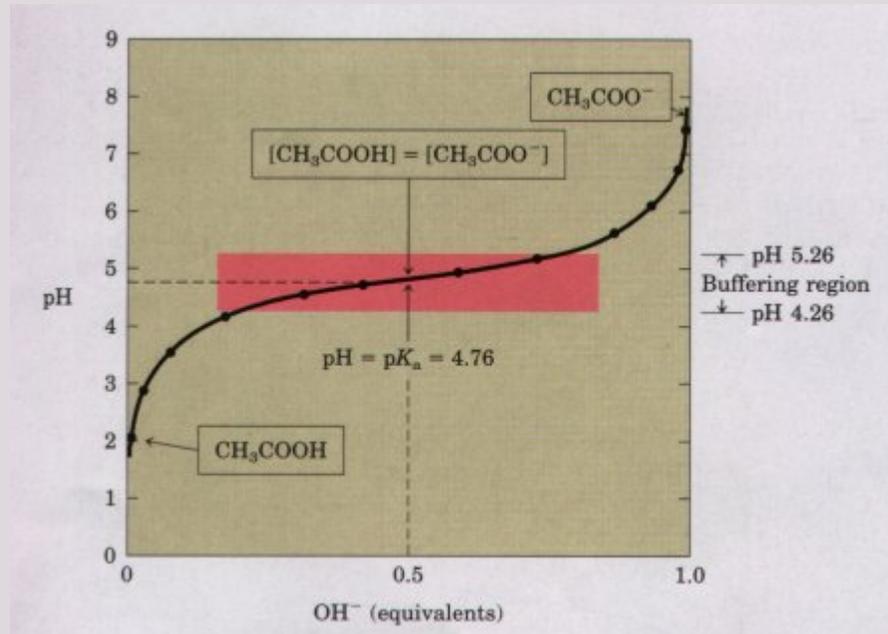
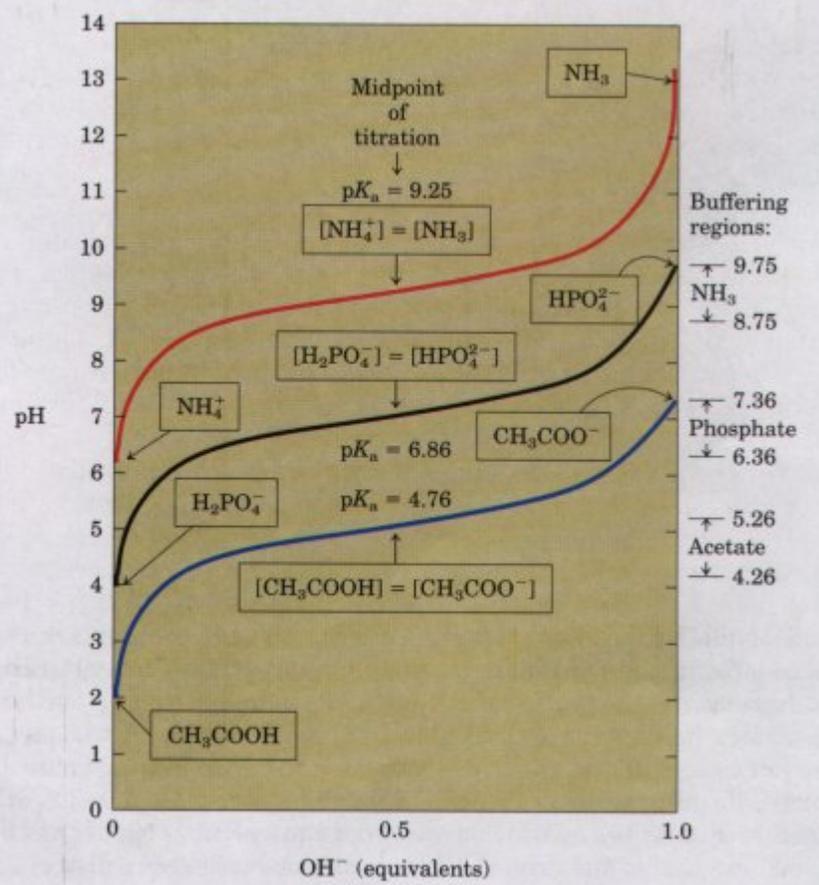
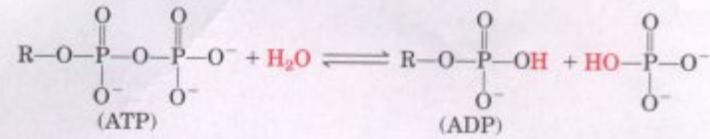


Table 4-7 Dissociation constant and pK_a of some common weak acids (proton donors) at 25 °C

Acid	K_a (M)	pK_a
HCOOH (formic acid)	1.78×10^{-4}	3.75
CH ₃ COOH (acetic acid)	1.74×10^{-5}	4.76
CH ₃ CH ₂ COOH (propionic acid)	1.35×10^{-5}	4.87
CH ₃ CH(OH)COOH (lactic acid)	1.38×10^{-4}	3.86
H ₃ PO ₄ (phosphoric acid)	7.25×10^{-3}	2.14
H ₂ PO ₄ ⁻ (dihydrogen phosphate)	1.38×10^{-7}	6.86
HPO ₄ ²⁻ (monohydrogen phosphate)	3.98×10^{-13}	12.4
H ₂ CO ₃ (carbonic acid)	1.70×10^{-4}	3.77
HCO ₃ ⁻ (bicarbonate)	6.31×10^{-11}	10.2
NH ₄ ⁺ (ammonium)	5.62×10^{-10}	9.25

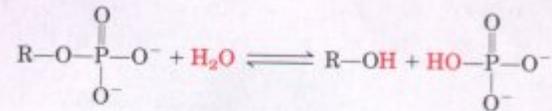






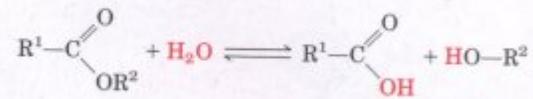
Phosphate anhydride

(a)



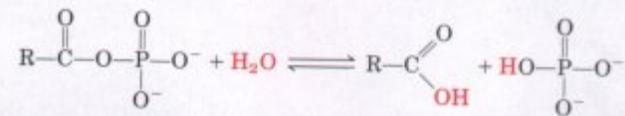
Phosphate ester

(b)



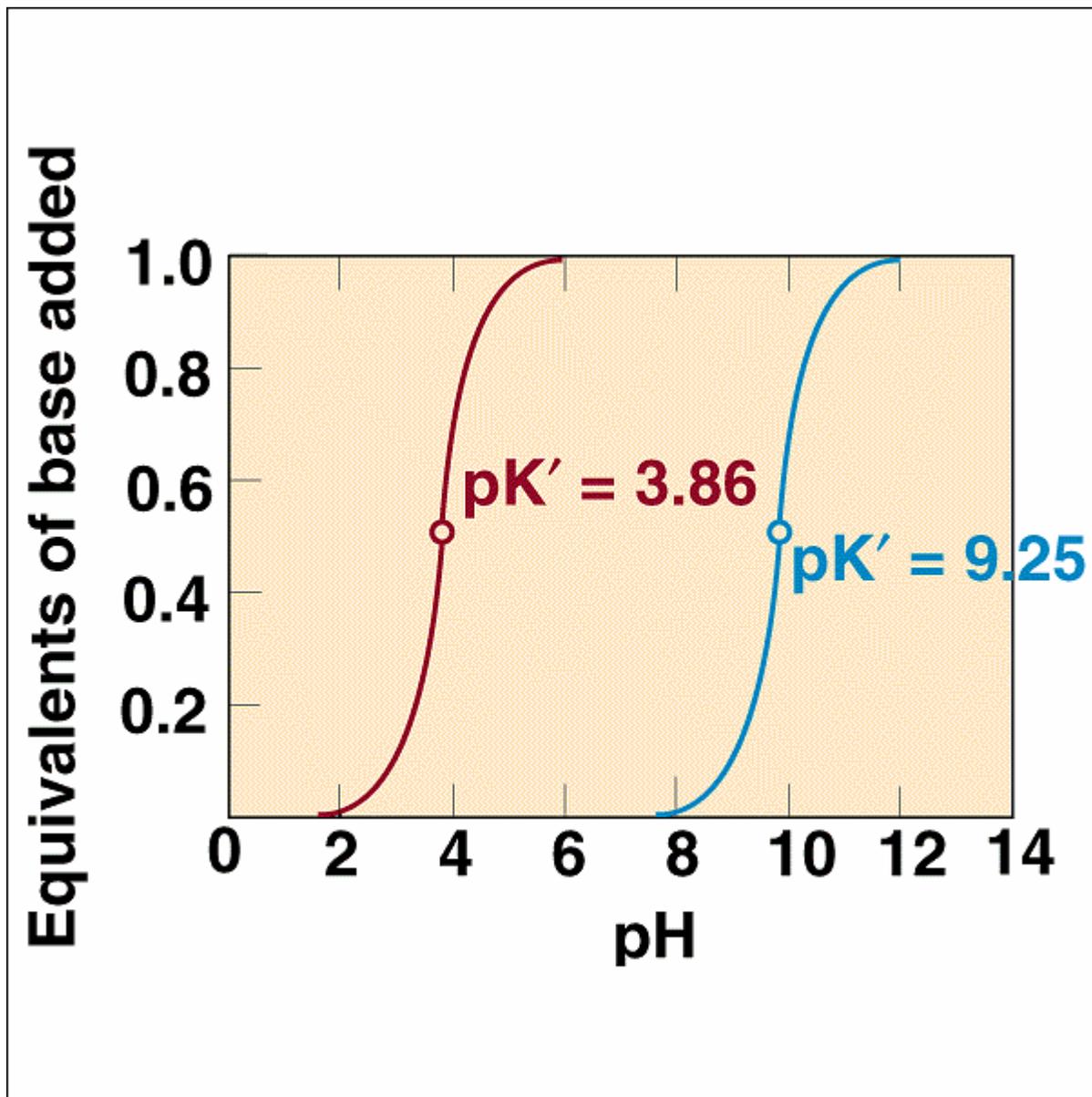
Carboxylate ester

(c)

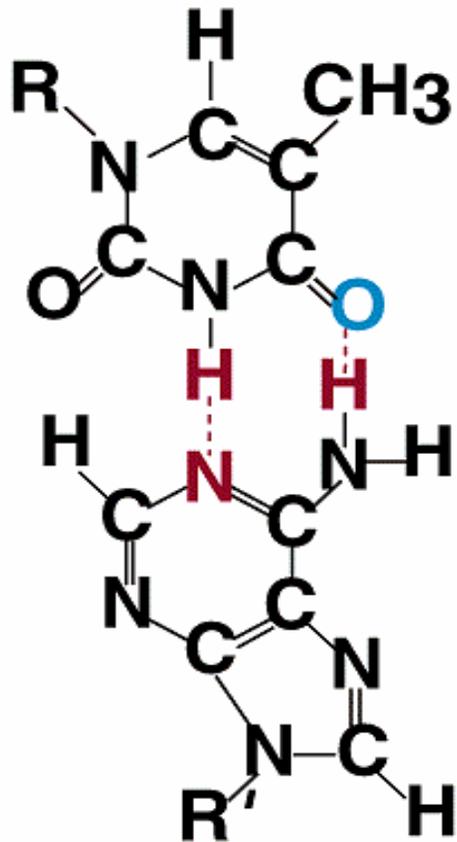


Acylphosphate anhydride

(d)







Thymine in DNA

Adenine in DNA

