## HEAT OF COMBUSTION

The heat of combustion of a substance at 25°C can be calculated from the enthalpy of formation  $(\Delta_t H^\circ)$  data in the table "Standard Thermodynamic Properties of Chemical Substances" in this Section. We can write the general combustion reaction as

$$X + O_2 \rightarrow CO_2(g) + H_2O(l) + other products$$

For a compound containing only carbon, hydrogen, and oxygen, the reaction is simply

$$C_a H_b O_c + \left(a + \frac{1}{4}b - \frac{1}{2}c\right) O_2 \rightarrow a CO_2(g) + \frac{1}{2}b H_2 O(l)$$

and the standard heat of combustion  $\Delta_c H^\circ$ , which is defined as the negative of the enthalpy change for the reaction (i.e., the heat released in the combustion process), is given by

| Molecular formula               | Name                | $\Delta_{c} H^{\circ}/kJ \text{ mol}^{-1}$ |
|---------------------------------|---------------------|--------------------------------------------|
| Inorganic substances            |                     |                                            |
| С                               | Carbon (graphite)   | 393.5                                      |
| CO                              | Carbon monoxide (g) | 283.0                                      |
| H <sub>2</sub>                  | Hydrogen (g)        | 285.8                                      |
| H <sub>3</sub> N                | Ammonia (g)         | 382.8                                      |
| $H_4N_2$                        | Hydrazine (g)       | 667.1                                      |
| N <sub>2</sub> O                | Nitrous oxide (g)   | 82.1                                       |
| Hydrocarbons                    |                     |                                            |
| $CH_4$                          | Methane (g)         | 890.8                                      |
| C <sub>2</sub> H <sub>2</sub>   | Acetylene (g)       | 1301.1                                     |
| $C_2H_4$                        | Ethylene (g)        | 1411.2                                     |
| $C_2H_6$                        | Ethane (g)          | 1560.7                                     |
| $C_{3}H_{6}$                    | Propylene (g)       | 2058.0                                     |
| $C_{3}H_{6}$                    | Cyclopropane (g)    | 2091.3                                     |
| $C_{3}H_{8}$                    | Propane (g)         | 2219.2                                     |
| $C_4H_6$                        | 1,3-Butadiene (g)   | 2541.5                                     |
| $C_4 H_{10}$                    | Butane (g)          | 2877.6                                     |
| $C_{5}H_{12}$                   | Pentane (l)         | 3509.0                                     |
| $C_6H_6$                        | Benzene (l)         | 3267.6                                     |
| $C_{6}H_{12}$                   | Cyclohexane (l)     | 3919.6                                     |
| $C_{6}H_{14}$                   | Hexane (l)          | 4163.2                                     |
| $C_7 H_8$                       | Toluene (l)         | 3910.3                                     |
| $C_{7}H_{16}$                   | Heptane (l)         | 4817.0                                     |
| $C_{10}H_{8}$                   | Naphthalene (s)     | 5156.3                                     |
| Alcohols and ethers             |                     |                                            |
| CH <sub>4</sub> O               | Methanol (l)        | 726.1                                      |
| C <sub>2</sub> H <sub>6</sub> O | Ethanol (l)         | 1366.8                                     |
| C,HO                            | Dimethyl ether (g)  | 1460.4                                     |
| C,HO,                           | Ethylene glycol (l) | 1189.2                                     |
| $C_3H_8O^2$                     | 1-Propanol (l)      | 2021.3                                     |

$$\Delta_c H^o = -a\Delta_f H^o(\mathrm{CO}_2, \mathbf{g}) - \frac{1}{2}b\Delta_f H^o(\mathrm{H}_2\mathrm{O}, \mathbf{l}) + \Delta_f H^o(\mathrm{C}_a\mathrm{H}_b\mathrm{O}_c)$$
  
= 393.51a + 142.915b +  $\Delta_f H^o(\mathrm{C}_a\mathrm{H}_b\mathrm{O}_c)$ 

This equation applies if the reactants start in their standard states (25°C and one atmosphere pressure) and the products return to the same conditions. The same equation applies to a compound containing another element if that element ends in its standard reference state (e.g., nitrogen, if the product is  $N_2$ ); in general, however, the exact products containing the other elements must be known in order to calculate the heat of combustion.

The following table gives the standard heat of combustion calculated in this manner for a few representative substances.

| Molecular formula                | Name                 | $\Delta_c H^{\circ}/kJ \text{ mol}^{-1}$ |
|----------------------------------|----------------------|------------------------------------------|
| $C_3H_8O_3$                      | Glycerol (l)         | 1655.4                                   |
| $C_4 H_{10} O$                   | Diethyl ether (l)    | 2723.9                                   |
| $C_5 H_{12} O$                   | 1-Pentanol (l)       | 3330.9                                   |
| C <sub>6</sub> H <sub>6</sub> O  | Phenol (s)           | 3053.5                                   |
| Carbonyl compounds               |                      |                                          |
| CH,O                             | Formaldehyde (g)     | 570.7                                    |
| C,H,O                            | Ketene (g)           | 1025.4                                   |
| C,H,O                            | Acetaldehyde (l)     | 1166.9                                   |
| C <sub>3</sub> H <sub>6</sub> O  | Acetone (l)          | 1789.9                                   |
| C <sub>3</sub> H <sub>6</sub> O  | Propanal (l)         | 1822.7                                   |
| $C_4H_8O$                        | 2-Butanone (l)       | 2444.1                                   |
| Acids and esters                 |                      |                                          |
| CH <sub>2</sub> O <sub>2</sub>   | Formic acid (l)      | 254.6                                    |
| $C_{2}H_{4}O_{2}$                | Acetic acid (l)      | 874.2                                    |
| $C_2H_4O_2$                      | Methyl formate (l)   | 972.6                                    |
| $C_3H_6O_2$                      | Methyl acetate (l)   | 1592.2                                   |
| $C_4H_8O_2$                      | Ethyl acetate (l)    | 2238.1                                   |
| $C_6H_5NO_2$                     | Nicotinic acid (s)   | 2731.1                                   |
| $C_7 H_6 O_2$                    | Benzoic acid (s)     | 3228.2                                   |
| Nitrogen compounds               |                      |                                          |
| CHN                              | Hydrogen cyanide (g) | 671.5                                    |
| CH <sub>3</sub> NO <sub>2</sub>  | Nitromethane (l)     | 709.2                                    |
| CH <sub>4</sub> N <sub>2</sub> O | Urea (s)             | 632.7                                    |
| CH₅N                             | Methylamine (g)      | 1085.6                                   |
| $C_2H_3N$                        | Acetonitrile (l)     | 1247.2                                   |
| $C_2H_5NO$                       | Acetamide (s)        | 1184.6                                   |
| $C_{3}H_{9}N$                    | Trimethylamine (g)   | 2443.1                                   |
| $C_5H_5N$                        | Pyridine (l)         | 2782.3                                   |
| C <sub>6</sub> H <sub>7</sub> N  | Aniline (l)          | 3392.8                                   |