

# 3-Methylenecyclopentene

|                      |                                            |
|----------------------|--------------------------------------------|
| Other names:         | Cyclopentene,3-methylene-                  |
| Inchi:               | InChI=1S/C6H8/c1-6-4-2-3-5-6/h2,4H,1,3,5H2 |
| InchiKey:            | YWQLRBQGXXHZJCF-UHFFFAOYSA-N               |
| Formula:             | C6H8                                       |
| SMILES:              | C=C1C=CCC1                                 |
| Mol. weight [g/mol]: | 80.13                                      |
| CAS:                 | 930-26-7                                   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 126.94  | kJ/mol  | Joback Method  |
| hf            | 115.00  | kJ/mol  | NIST Webbook   |
| hfus          | 4.22    | kJ/mol  | Joback Method  |
| hvap          | 29.97   | kJ/mol  | Joback Method  |
| ie            | 8.40    | eV      | NIST Webbook   |
| log10ws       | -1.94   |         | Crippen Method |
| logp          | 1.893   |         | Crippen Method |
| mcvol         | 75.940  | ml/mol  | McGowan Method |
| pc            | 4294.29 | kPa     | Joback Method  |
| tb            | 354.95  | K       | Joback Method  |
| tc            | 554.62  | K       | Joback Method  |
| tf            | 186.96  | K       | Joback Method  |
| vc            | 0.283   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 116.01 | J/mol×K | 354.95          | Joback Method |
| cpg           | 163.49 | J/mol×K | 521.34          | Joback Method |
| cpg           | 155.05 | J/mol×K | 488.06          | Joback Method |
| cpg           | 146.11 | J/mol×K | 454.79          | Joback Method |
| cpg           | 136.64 | J/mol×K | 421.51          | Joback Method |
| cpg           | 126.62 | J/mol×K | 388.23          | Joback Method |
| cpg           | 171.43 | J/mol×K | 554.62          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002661 | Pa×s | 354.95 | Joback Method |
| dvisc | 0.0003217 | Pa×s | 326.95 | Joback Method |
| dvisc | 0.0004029 | Pa×s | 298.95 | Joback Method |
| dvisc | 0.0005286 | Pa×s | 270.95 | Joback Method |
| dvisc | 0.0007384 | Pa×s | 242.96 | Joback Method |
| dvisc | 0.0011253 | Pa×s | 214.96 | Joback Method |
| dvisc | 0.0019455 | Pa×s | 186.96 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C930267&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C930267&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                 |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>ie:</b>      | Ionization energy                               |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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