

# Cholest-5-en-3-one

|                             |   |
|-----------------------------|---|
| <b>Other names:</b>         | Cholesterone<br>Oxidized cholesterol<br>5-Cholesten-3-one<br>«delta»(sup5)-Cholestenone<br>Cholestenone |
| <b>Inchi:</b>               | InChI=1S/C27H44O/c1-18(2)7-6-8-19(3)23-11-12-24-22-10-9-20-17-21(28)13-15-26(20,4                       |
| <b>InchiKey:</b>            | GGCLNOIGPMGLDB-XUJZTTFYSA-N   |
| <b>Formula:</b>             | C27H44O   |
| <b>SMILES:</b>              | CC(C)CCCC(C)C1CCC2C3CC=C4CC(=O)CCC4(C)C3CCC12C  |
| <b>Mol. weight [g/mol]:</b> | 384.64  |
| <b>CAS:</b>                 | 601-54-7  |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 225.42  | kJ/mol  | Joback Method  |
| hf            | -452.36 | kJ/mol  | Joback Method  |
| hfus          | 30.57   | kJ/mol  | Joback Method  |
| hvap          | 77.71   | kJ/mol  | Joback Method  |
| log10ws       | -7.90   |         | Crippen Method |
| logp          | 7.597   |         | Crippen Method |
| mcvol         | 345.120 | ml/mol  | McGowan Method |
| pc            | 1061.71 | kPa     | Joback Method  |
| tb            | 927.69  | K       | Joback Method  |
| tc            | 1161.44 | K       | Joback Method  |
| tf            | 539.03  | K       | Joback Method  |
| vc            | 1.310   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value   | Unit    | Temperature [K] | Source        |
|---------------|---------|---------|-----------------|---------------|
| cpg           | 1275.88 | J/molxK | 927.69          | Joback Method |
| cpg           | 1307.11 | J/molxK | 966.65          | Joback Method |
| cpg           | 1338.52 | J/molxK | 1005.61         | Joback Method |
| cpg           | 1370.45 | J/molxK | 1044.56         | Joback Method |

|     |         |         |         |               |
|-----|---------|---------|---------|---------------|
| cpg | 1403.26 | J/mol×K | 1083.52 | Joback Method |
| cpg | 1437.28 | J/mol×K | 1122.48 | Joback Method |
| cpg | 1472.84 | J/mol×K | 1161.44 | 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=C601547&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C601547&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>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>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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