

# Cedrenol acetate

**Other names:**

**Inchi:** [3R,1(3«alpha»,3a«beta»,7«beta»,8a«alpha»)-2,3,4,7,8,8a-hexahydro-3,8,8-trimethyl-1H-  
acetal]C17H26O2/c1-10-6-7-15-16(4,5)13-8-17(10,15)9-14(11(13)2)19-12(3)18/10,1

InchiKey: UTZBRSREMTWJBB-OOXPIDESA-N

**Formula:** C<sub>17</sub>H<sub>26</sub>O<sub>2</sub>

**SMILES:** C=C1C(OC(C)=O)CC23CC1C(C)(C)C2CCC3C

Mol. weight [g/mol]: 262.39

**CAS:** 1405-92-1

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 35.36   | kJ/mol  | Joback Method  |
| hf            | -379.23 | kJ/mol  | Joback Method  |
| hfus          | 22.24   | kJ/mol  | Joback Method  |
| hvap          | 59.60   | kJ/mol  | Joback Method  |
| log10ws       | -4.24   |         | Crippen Method |
| logp          | 3.957   |         | Crippen Method |
| mcvol         | 220.950 | ml/mol  | McGowan Method |
| pc            | 1804.63 | kPa     | Joback Method  |
| rinpol        | 1743.00 |         | NIST Webbook   |
| rinpol        | 1743.00 |         | NIST Webbook   |
| tb            | 679.04  | K       | Joback Method  |
| tc            | 899.33  | K       | Joback Method  |
| tf            | 449.05  | K       | Joback Method  |
| vc            | 0.844   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 669.85 | J/mol×K | 679.04          | Joback Method |
| cpg           | 691.60 | J/mol×K | 715.75          | Joback Method |
| cpg           | 712.47 | J/mol×K | 752.47          | Joback Method |
| cpg           | 732.71 | J/mol×K | 789.18          | Joback Method |
| cpg           | 752.57 | J/mol×K | 825.90          | Joback Method |
| cpg           | 772.32 | J/mol×K | 862.61          | 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=C1405921&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1405921&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>rinpol:</b>  | Non-polar retention indices                     |
| <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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