

# Caryophylle-4(12),8(13)-dien-5«beta»-ol

|                      |                                                                                   |
|----------------------|-----------------------------------------------------------------------------------|
| Inchi:               | InChI=1S/C15H24O/c1-10-6-8-14(16)11(2)5-7-13-12(10)9-15(13,3)4/h12-14,16H,1-2,5-9 |
| InchiKey:            | CIIYOYPOMGIECX-MELADBBJSA-N                                                       |
| Formula:             | C15H24O                                                                           |
| SMILES:              | C=C1CCC2C(CC2(C)C)C(=C)CCC1O                                                      |
| Mol. weight [g/mol]: | 220.35                                                                            |

## Physical Properties

| Property code | Value   | Unit   | Source             |
|---------------|---------|--------|--------------------|
| hfus          | 17.99   | kJ/mol | Joback Method      |
| hvap          | 91.70   | kJ/mol | Cheméo Relay (1.0) |
| logp          | 3.696   |        | Crippen Method     |
| mcvol         | 197.760 | ml/mol | McGowan Method     |
| rinpol        | 1641.00 |        | NIST Webbook       |
| rinpol        | 1590.00 |        | NIST Webbook       |
| rinpol        | 1641.00 |        | NIST Webbook       |
| tb            | 577.46  | K      | Cheméo Relay (1.0) |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 575.21 | J/mol×K | 658.83          | Joback Method |
| cpg           | 594.89 | J/mol×K | 693.49          | Joback Method |
| cpg           | 613.57 | J/mol×K | 728.16          | Joback Method |
| cpg           | 631.35 | J/mol×K | 762.82          | Joback Method |
| cpg           | 648.34 | J/mol×K | 797.48          | Joback Method |
| cpg           | 664.64 | J/mol×K | 832.14          | Joback Method |
| cpg           | 680.34 | J/mol×K | 866.81          | Joback Method |

## Sources

Crippen Method: [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

**Cheméo Relay (1.0):**

**Cheméo Relay (1.0, beta):**

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=R612807&Units=SI>

**Joback Method:** [https://en.wikipedia.org/wiki/Joback\\_method](https://en.wikipedia.org/wiki/Joback_method)

**McGowan Method:** <http://link.springer.com/article/10.1007/BF02311772>

**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci990307l>

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>rinpola:</b> | Non-polar retention indices                     |
| <b>tb:</b>      | Normal Boiling Point Temperature                |

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