

(E)-Caryophyllene epoxide

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| Other names: | Caryophyllene oxide 2 |
| Inchi: | InChI=1S/C15H24O/c1-10-5-6-13-15(4,16-13)8-7-12-11(10)9-14(12,2)3/h11-13H,1,5-9H2 |
| InchiKey: | NVEQFIOZRFFVFW-YKQCJFHZSA-N |
| Formula: | C15H24O |
| SMILES: | <chem>C=C1CCC2OC2(C)CCC2C1CC2(C)C</chem> |
| Mol. weight [g/mol]: | 220.35 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 161.93 | kJ/mol | Joback Method |
| hf | -210.97 | kJ/mol | Joback Method |
| hfus | 19.08 | kJ/mol | Joback Method |
| hvap | 50.99 | kJ/mol | Joback Method |
| log10ws | -4.23 | | Crippen Method |
| logp | 3.936 | | Crippen Method |
| mcvol | 191.200 | ml/mol | McGowan Method |
| pc | 2145.33 | kPa | Joback Method |
| rinpol | 1564.00 | | NIST Webbook |
| rinpol | 1570.00 | | NIST Webbook |
| ripol | 2010.00 | | NIST Webbook |
| ripol | 2010.00 | | NIST Webbook |
| ripol | 1960.00 | | NIST Webbook |
| ripol | 2019.00 | | NIST Webbook |
| ripol | 1976.00 | | NIST Webbook |
| tb | 592.88 | K | Joback Method |
| tc | 822.50 | K | Joback Method |
| tf | 381.64 | K | Joback Method |
| vc | 0.722 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 539.37 | J/molxK | 592.88 | Joback Method |
| cpg | 562.46 | J/molxK | 631.15 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 584.07 | J/mol×K | 669.42 | Joback Method |
| cpg | 604.52 | J/mol×K | 707.69 | Joback Method |
| cpg | 624.09 | J/mol×K | 745.96 | Joback Method |
| cpg | 643.08 | J/mol×K | 784.23 | Joback Method |
| cpg | 661.78 | J/mol×K | 822.50 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R328002&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpola: | Non-polar retention indices |
| ripola: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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<https://www.chemeo.com/cid/75-678-7/E-Caryophyllene-epoxide.pdf>

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