

6-epi-6,9-Epoxyfarnesa-1,7(14),10-trien-3-ol

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|-----------------------------|--|
| Inchi: | InChI=1S/C15H24O2/c1-6-15(5,16)8-7-14-12(4)10-13(17-14)9-11(2)3/h6,9,13-14,16H,1,4 |
| InchiKey: | LJPFBTWRZZGKFP-UHFFFAOYSA-N |
| Formula: | C15H24O2 |
| SMILES: | C=CC(C)(O)CCC1OC(C=C(C)C)CC1=C |
| Mol. weight [g/mol]: | 236.35 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 96.75 | kJ/mol | Joback Method |
| hf | -288.67 | kJ/mol | Joback Method |
| hfus | 30.72 | kJ/mol | Joback Method |
| hvap | 68.35 | kJ/mol | Joback Method |
| log10ws | -4.24 | | Crippen Method |
| logp | 3.383 | | Crippen Method |
| mcvol | 210.190 | ml/mol | McGowan Method |
| pc | 1905.24 | kPa | Joback Method |
| rinpol | 1614.00 | | NIST Webbook |
| rinpol | 1614.00 | | NIST Webbook |
| rinpol | 1614.00 | | NIST Webbook |
| ripol | 2284.00 | | NIST Webbook |
| ripol | 2276.00 | | NIST Webbook |
| tb | 668.99 | K | Joback Method |
| tc | 862.55 | K | Joback Method |
| tf | 348.16 | K | Joback Method |
| vc | 0.790 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 598.85 | J/molxK | 668.99 | Joback Method |
| cpg | 615.48 | J/molxK | 701.25 | Joback Method |
| cpg | 631.15 | J/molxK | 733.51 | Joback Method |
| cpg | 645.92 | J/molxK | 765.77 | Joback Method |
| cpg | 659.85 | J/molxK | 798.03 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 672.99 | J/mol×K | 830.29 | Joback Method |
| cpg | 685.41 | J/mol×K | 862.55 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R232525&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |

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 |
| hvp: | 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 |
| rinpol: | Non-polar retention indices |
| ripol: | 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/41-668-5/6-epi-6-9-Epoxyfarnesa-1-7-14-10-trien-3-ol.pdf>

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