

«epsilon»-Bulgarene

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|-----------------------------|---|
| Inchi: | InChI=1S/C15H24/c1-10(2)13-8-6-12(4)14-7-5-11(3)9-15(13)14/h10,13-15H,3-9H2,1-2H3 |
| InchiKey: | NOLWRMQDWRAODO-ZNMIVQPWSA-N |
| Formula: | C15H24 |
| SMILES: | C=C1CCC2C(=C)CCC(C(C)C)C2C1 |
| Mol. weight [g/mol]: | 204.35 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 244.53 | kJ/mol | Joback Method |
| hf | -89.11 | kJ/mol | Joback Method |
| hfus | 17.71 | kJ/mol | Joback Method |
| hvap | 49.12 | kJ/mol | Joback Method |
| log10ws | -4.63 | | Crippen Method |
| logp | 4.581 | | Crippen Method |
| mcvol | 191.890 | ml/mol | McGowan Method |
| pc | 1915.26 | kPa | Joback Method |
| rinpol | 1514.00 | | NIST Webbook |
| rinpol | 1514.00 | | NIST Webbook |
| tb | 566.37 | K | Joback Method |
| tc | 779.32 | K | Joback Method |
| tf | 288.73 | K | Joback Method |
| vc | 0.719 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 499.92 | J/molxK | 566.37 | Joback Method |
| cpg | 603.18 | J/molxK | 743.83 | Joback Method |
| cpg | 584.99 | J/molxK | 708.34 | Joback Method |
| cpg | 565.61 | J/molxK | 672.85 | Joback Method |
| cpg | 545.00 | J/molxK | 637.35 | Joback Method |
| cpg | 523.12 | J/molxK | 601.86 | Joback Method |
| cpg | 620.22 | J/molxK | 779.32 | Joback Method |
| dvisc | 0.0004074 | Paxs | 566.37 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0004790 | Paxs | 520.10 | Joback Method |
| dvisc | 0.0005812 | Paxs | 473.82 | Joback Method |
| dvisc | 0.0007353 | Paxs | 427.55 | Joback Method |
| dvisc | 0.0009850 | Paxs | 381.28 | Joback Method |
| dvisc | 0.0014305 | Paxs | 335.00 | Joback Method |
| dvisc | 0.0023414 | Paxs | 288.73 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=R206947&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

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|----------------------------|---|
| cp_g: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| g_f: | Standard Gibbs free energy of formation |
| h_f: | Enthalpy of formation at standard conditions |
| h_{fus}: | Enthalpy of fusion at standard conditions |
| h_{vap}: | Enthalpy of vaporization at standard conditions |
| log₁₀ws: | Log ₁₀ of Water solubility in mol/l |
| log_p: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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