

Germacranolide callitris

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|-----------------------------|---|
| Inchi: | InChI=1S/C15H22O2/c1-10-5-4-6-11(2)9-14-13(8-7-10)12(3)15(16)17-14/h6-7,12-14H,4- |
| InchiKey: | QJRFOUJEGHRZIU-NXAIOARDSA-N |
| Formula: | C15H22O2 |
| SMILES: | CC1=CCC2C(CC(C)=CCC1)OC(=O)C2C |
| Mol. weight [g/mol]: | 234.33 |
| CAS: | 57759-35-0 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -63.54 | kJ/mol | Joback Method |
| hf | -447.87 | kJ/mol | Joback Method |
| hfus | 26.40 | kJ/mol | Joback Method |
| hvap | 60.37 | kJ/mol | Joback Method |
| log10ws | -4.09 | | Crippen Method |
| logp | 3.631 | | Crippen Method |
| mcvol | 199.330 | ml/mol | McGowan Method |
| pc | 2088.84 | kPa | Joback Method |
| rinpol | 1972.80 | | NIST Webbook |
| rinpol | 1972.80 | | NIST Webbook |
| tb | 684.35 | K | Joback Method |
| tc | 927.07 | K | Joback Method |
| tf | 387.16 | K | Joback Method |
| vc | 0.733 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 586.87 | J/mol×K | 684.35 | Joback Method |
| cpg | 609.93 | J/mol×K | 724.80 | Joback Method |
| cpg | 631.32 | J/mol×K | 765.26 | Joback Method |
| cpg | 651.02 | J/mol×K | 805.71 | Joback Method |
| cpg | 669.03 | J/mol×K | 846.17 | Joback Method |
| cpg | 685.32 | J/mol×K | 886.62 | Joback Method |
| cpg | 699.89 | J/mol×K | 927.07 | Joback Method |

Sources

| | |
|------------------------|---|
| 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 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C57759350&Units=SI |

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 |
| rinp: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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