

Sesquisabinene hydrate

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|-----------------------------|---|
| Inchi: | InChI=1S/C15H26O/c1-11(2)6-5-7-12(3)15-9-8-14(4,16)13(15)10-15/h6,12-13,16H,5,7-1 |
| InchiKey: | IRDFGGRWKUKANK-NEXFUWMNSA-N |
| Formula: | C15H26O |
| SMILES: | CC(C)=CCCC(C)C12CCC(C)(O)C1C2 |
| Mol. weight [g/mol]: | 222.37 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 110.64 | kJ/mol | Joback Method |
| hf | -247.27 | kJ/mol | Joback Method |
| hfus | 18.81 | kJ/mol | Joback Method |
| hvap | 62.53 | kJ/mol | Joback Method |
| log10ws | -4.39 | | Crippen Method |
| logp | 3.920 | | Crippen Method |
| mcvol | 202.060 | ml/mol | McGowan Method |
| pc | 2096.50 | kPa | Joback Method |
| rinpol | 1579.00 | | NIST Webbook |
| rinpol | 1588.00 | | NIST Webbook |
| rinpol | 1605.00 | | NIST Webbook |
| rinpol | 1544.00 | | NIST Webbook |
| rinpol | 1579.00 | | NIST Webbook |
| rinpol | 1547.00 | | NIST Webbook |
| rinpol | 1547.00 | | NIST Webbook |
| ripol | 2018.00 | | NIST Webbook |
| ripol | 1997.00 | | NIST Webbook |
| tb | 647.67 | K | Joback Method |
| tc | 844.14 | K | Joback Method |
| tf | 365.03 | K | Joback Method |
| vc | 0.778 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 583.77 | J/mol×K | 647.67 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 600.91 | J/mol×K | 680.41 | Joback Method |
| cpg | 617.33 | J/mol×K | 713.16 | Joback Method |
| cpg | 633.24 | J/mol×K | 745.90 | Joback Method |
| cpg | 648.84 | J/mol×K | 778.65 | Joback Method |
| cpg | 664.35 | J/mol×K | 811.39 | Joback Method |
| cpg | 679.97 | J/mol×K | 844.14 | 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=R313976&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

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 |
| ripol: | 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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