

Gymnomitr-3-en-15-ol

| | |
|-----------------------------|---|
| Inchi: | InChI=1S/C15H24O/c1-13-8-5-11(10-16)12(9-13)14(2)6-4-7-15(13,14)3/h5,12,16H,4,6-1 |
| InchiKey: | BOBUWMAINZVWRJ-UHFFFAOYSA-N |
| Formula: | C15H24O |
| SMILES: | CC12CC=C(CO)C(C1)C1(C)CCCC21C |
| Mol. weight [g/mol]: | 220.35 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 92.80 | kJ/mol | Joback Method |
| hf | -227.39 | kJ/mol | Joback Method |
| hfus | 11.91 | kJ/mol | Joback Method |
| hvap | 62.94 | kJ/mol | Joback Method |
| log10ws | -3.94 | | Crippen Method |
| logp | 3.531 | | Crippen Method |
| mcvol | 191.200 | ml/mol | McGowan Method |
| pc | 2492.52 | kPa | Joback Method |
| rinqol | 1719.00 | | NIST Webbook |
| tb | 663.73 | K | Joback Method |
| tc | 879.86 | K | Joback Method |
| tf | 447.15 | K | Joback Method |
| vc | 0.729 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 568.91 | J/mol×K | 663.73 | Joback Method |
| cpg | 586.84 | J/mol×K | 699.75 | Joback Method |
| cpg | 604.22 | J/mol×K | 735.77 | Joback Method |
| cpg | 621.43 | J/mol×K | 771.80 | Joback Method |
| cpg | 638.80 | J/mol×K | 807.82 | Joback Method |
| cpg | 656.69 | J/mol×K | 843.84 | Joback Method |
| cpg | 675.47 | J/mol×K | 879.86 | Joback Method |

Sources

| | |
|------------------------|---|
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| 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=R429546&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 |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mccvol: | 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 |

Latest version available from:

<https://www.chemeo.com/cid/14-109-5/Gymnomitri-3-en-15-ol.pdf>

Generated by Cheméo on 2024-04-19 01:31:34.374792641 +0000 UTC m=+15779543.295369956.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.