

(1R,1'R

)-1-(6,6-dimethyl-bicyclo[3.1.1]hept-2-en-

2-yl)-1-hydroxy-2-propanone, R

Formula:

C₁₂H₁₈O₂

SMILES:

CC(=O)C(O)C1=CCC2CC1C2(C)C

Mol. weight [g/mol]:

194.27

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -101.49 | kJ/mol | Joback Method |
| hf | -380.45 | kJ/mol | Joback Method |
| hfus | 18.78 | kJ/mol | Joback Method |
| hvap | 64.83 | kJ/mol | Joback Method |
| log10ws | -2.42 | | Crippen Method |
| logp | 1.929 | | Crippen Method |
| mcvol | 161.360 | ml/mol | McGowan Method |
| pc | 2802.44 | kPa | Joback Method |
| rinpol | 1362.00 | | NIST Webbook |
| rinpol | 1362.00 | | NIST Webbook |
| ripol | 1966.00 | | NIST Webbook |
| tb | 637.03 | K | Joback Method |
| tc | 839.39 | K | Joback Method |
| tf | 386.05 | K | Joback Method |
| vc | 0.616 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 452.88 | J/mol×K | 637.03 | Joback Method |
| cpg | 467.24 | J/mol×K | 670.76 | Joback Method |
| cpg | 480.86 | J/mol×K | 704.48 | Joback Method |
| cpg | 493.86 | J/mol×K | 738.21 | Joback Method |
| cpg | 506.35 | J/mol×K | 771.94 | Joback Method |
| cpg | 518.48 | J/mol×K | 805.67 | Joback Method |
| cpg | 530.37 | J/mol×K | 839.39 | 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=R522600&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 |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpolar: | Non-polar retention indices |
| ripolar: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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