

Cyclopentanol, 1,2-dimethyl-3-(1-methylethenyl)-, [1R-(1«alpha»,2«beta»,3«beta»)]-

Other names: Cyclopentanol, 3-isopropenyl-1,1-dimethyl-, (1R,2R,3R)-(+)-
Plinol D

Inchi: Plinol D, (+)-
InChI=1S/C10H18O/c1-7(2)9-5-6-10(4,11)8(9)3/h8-9,11H,1,5-6H2,2-4H3

InchiKey: ZRVPDCMGGOSDKG-UHFFFAOYSA-N

Formula: C10H18O

SMILES: C=C(C)C1CCC(C)(O)C1C

Mol. weight [g/mol]: 154.25

CAS: 4028-58-4

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | -8.57 | kJ/mol | Joback Method |
| hf | -251.28 | kJ/mol | Joback Method |
| hfus | 12.93 | kJ/mol | Joback Method |
| hvap | 52.43 | kJ/mol | Joback Method |
| log10ws | -2.65 | | Crippen Method |
| logp | 2.360 | | Crippen Method |
| mcvol | 142.470 | ml/mol | McGowan Method |
| pc | 2829.33 | kPa | Joback Method |
| rinpol | 1503.00 | | NIST Webbook |
| tb | 523.12 | K | Joback Method |
| tc | 716.53 | K | Joback Method |
| tf | 273.88 | K | Joback Method |
| vc | 0.533 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 349.90 | J/molxK | 523.12 | Joback Method |
| cpg | 365.68 | J/molxK | 555.35 | Joback Method |
| cpg | 380.58 | J/molxK | 587.59 | Joback Method |
| cpg | 394.67 | J/molxK | 619.82 | Joback Method |
| cpg | 408.03 | J/molxK | 652.06 | Joback Method |

| | | | | |
|-----|--------|---------|--------|---------------|
| cpg | 420.76 | J/mol×K | 684.29 | Joback Method |
| cpg | 432.93 | J/mol×K | 716.53 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C4028584&Units=SI |
| 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 |

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 |
| 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/71-655-6/Cyclopentanol-1-2-dimethyl-3-1-methylethenyl-1R-1-alpha-2-beta-3-beta.pdf>

Generated by Cheméo on 2024-04-27 19:24:27.062848676 +0000 UTC m=+16535115.983425987.

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