

# cis,cis,trans-1,2,4-Trimethylcyclopentane

|                      |   |
|----------------------|---|
| Inchi:               | InChI=1S/C8H16/c1-6-4-7(2)8(3)5-6/h6-8H,4-5H2,1-3H3/t6-,7+,8- |
| InchiKey:            | PNUFYSGVPVMNRN-RNLVFQAGSA-N                                   |
| Formula:             | C8H16   |
| SMILES:              | CC1CC(C)C(C)C1  |
| Mol. weight [g/mol]: | 112.21  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 37.61   | kJ/mol               | Joback Method  |
| hf            | -188.65 | kJ/mol               | Joback Method  |
| hfus          | 12.55   | kJ/mol               | Joback Method  |
| hvap          | 33.04   | kJ/mol               | Joback Method  |
| log10ws       | -2.34   |                      | Crippen Method |
| logp          | 2.688   |                      | Crippen Method |
| mvol          | 112.720 | ml/mol               | McGowan Method |
| pc            | 2853.57 | kPa                  | Joback Method  |
| rinpol        | 768.00  |                      | NIST Webbook   |
| rinpol        | 768.00  |                      | NIST Webbook   |
| tb            | 388.38  | K                    | Joback Method  |
| tc            | 579.97  | K                    | Joback Method  |
| tf            | 182.34  | K                    | Joback Method  |
| vc            | 0.422   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 213.62    | J/molxK | 388.38          | Joback Method |
| cpg           | 230.52    | J/molxK | 420.31          | Joback Method |
| cpg           | 246.72    | J/molxK | 452.24          | Joback Method |
| cpg           | 262.24    | J/molxK | 484.17          | Joback Method |
| cpg           | 277.08    | J/molxK | 516.11          | Joback Method |
| cpg           | 291.26    | J/molxK | 548.04          | Joback Method |
| cpg           | 304.79    | J/molxK | 579.97          | Joback Method |
| dvisc         | 0.0010560 | Paxs    | 182.34          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0006962 | Paxs | 216.68 | Joback Method |
| dvisc | 0.0005145 | Paxs | 251.02 | Joback Method |
| dvisc | 0.0004089 | Paxs | 285.36 | Joback Method |
| dvisc | 0.0003414 | Paxs | 319.70 | Joback Method |
| dvisc | 0.0002952 | Paxs | 354.04 | Joback Method |
| dvisc | 0.0002619 | Paxs | 388.38 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R634090&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R634090&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>                                 |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                     |

## Legend

|                                       |   |
|---------------------------------------|---|
| <b>cp<sub>g</sub>:</b>                | Ideal gas heat capacity                         |
| <b>dvisc:</b>                         | Dynamic viscosity                               |
| <b>g<sub>f</sub>:</b>                 | Standard Gibbs free energy of formation         |
| <b>h<sub>f</sub>:</b>                 | Enthalpy of formation at standard conditions    |
| <b>h<sub>fus</sub>:</b>               | Enthalpy of fusion at standard conditions       |
| <b>h<sub>vap</sub>:</b>               | Enthalpy of vaporization at standard conditions |
| <b>log<sub>10</sub>w<sub>s</sub>:</b> | Log <sub>10</sub> of Water solubility in mol/l  |
| <b>log<sub>p</sub>:</b>               | Octanol/Water partition coefficient             |
| <b>mcvol:</b>                         | McGowan's characteristic volume                 |
| <b>pc:</b>                            | Critical Pressure                               |
| <b>rinpol:</b>                        | Non-polar retention indices                     |
| <b>tb:</b>                            | Normal Boiling Point Temperature                |
| <b>tc:</b>                            | Critical Temperature                            |
| <b>tf:</b>                            | Normal melting (fusion) point                   |
| <b>vc:</b>                            | Critical Volume                                 |

Latest version available from:

<https://www.chemeo.com/cid/77-950-2/cis-cis-trans-1-2-4-Trimethylcyclopentane.pdf>

Generated by Cheméo on 2024-04-19 20:14:52.534970049 +0000 UTC m=+15846941.455547366.

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