

Tetrahydro-alpha--tri-cyclopentadiene

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|-----------------------------|--|
| Inchi: | InChI=1S/C15H22/c1-2-10-11(3-1)13-7-12(10)14-8-4-5-9(6-8)15(13)14/h8-15H,1-7H2 |
| InchiKey: | SUUFUDGKCFRARM-UHFFFAOYSA-N |
| Formula: | C15H22 |
| SMILES: | C1CC2C(C1)C1CC2C2C3CCC(C3)C12 |
| Mol. weight [g/mol]: | 202.34 |
| CAS: | 75172-86-0 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|----------|----------------------|----------------|
| chs | -8935.40 | kJ/mol | NIST Webbook |
| gf | 356.04 | kJ/mol | Joback Method |
| hf | -49.95 | kJ/mol | Joback Method |
| hfus | 28.49 | kJ/mol | Joback Method |
| hvap | 47.62 | kJ/mol | Joback Method |
| log10ws | -3.64 | | Crippen Method |
| logp | 3.715 | | Crippen Method |
| mcvol | 167.910 | ml/mol | McGowan Method |
| pc | 2252.53 | kPa | Joback Method |
| tb | 562.29 | K | Joback Method |
| tc | 783.52 | K | Joback Method |
| tf | 335.79 | K | Joback Method |
| vc | 0.657 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 495.75 | J/molxK | 562.29 | Joback Method |
| cpg | 520.98 | J/molxK | 599.16 | Joback Method |
| cpg | 544.34 | J/molxK | 636.03 | Joback Method |
| cpg | 566.01 | J/molxK | 672.91 | Joback Method |
| cpg | 586.16 | J/molxK | 709.78 | Joback Method |
| cpg | 604.97 | J/molxK | 746.65 | Joback Method |
| cpg | 622.62 | J/molxK | 783.52 | Joback Method |
| dvisc | 0.0019631 | Paxs | 335.79 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0031769 | Paxs | 373.54 | Joback Method |
| dvisc | 0.0047063 | Paxs | 411.29 | Joback Method |
| dvisc | 0.0065262 | Paxs | 449.04 | Joback Method |
| dvisc | 0.0086024 | Paxs | 486.79 | Joback Method |
| dvisc | 0.0108972 | Paxs | 524.54 | Joback Method |
| dvisc | 0.0133727 | Paxs | 562.29 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C75172860&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 |

Legend

| | |
|-----------------|---|
| chs: | Standard solid enthalpy of combustion |
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| 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 |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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