

Cyclopropane, 1,1'-ethenylidenebis-

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|-----------------------------|--|
| Other names: | Ethylene, 1,1-dicyclopropyl- 1,1-Dicyclopropylethene 1,1-Dicyclopropylethylene 1,1'-vinylidenebiscyclopropane |
| Inchi: | InChI=1S/C8H12/c1-6(7-2-3-7)8-4-5-8/h7-8H,1-5H2 |
| InchiKey: | RXCSODBXHNSSCQ-UHFFFAOYSA-N |
| Formula: | C8H12 |
| SMILES: | C=C(C1CC1)C1CC1 |
| Mol. weight [g/mol]: | 108.18 |
| CAS: | 822-93-5 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| affp | 904.70 | kJ/mol | NIST Webbook |
| basg | 875.80 | kJ/mol | NIST Webbook |
| gf | 217.27 | kJ/mol | Joback Method |
| hf | 52.79 | kJ/mol | Joback Method |
| hfus | 10.16 | kJ/mol | Joback Method |
| hvap | 32.64 | kJ/mol | Joback Method |
| ie | 8.87 | eV | NIST Webbook |
| ie | 8.08 | eV | NIST Webbook |
| log10ws | -2.33 | | Crippen Method |
| logp | 2.363 | | Crippen Method |
| mcvol | 97.560 | ml/mol | McGowan Method |
| pc | 3547.31 | kPa | Joback Method |
| tb | 392.48 | K | Joback Method |
| tc | 594.03 | K | Joback Method |
| tf | 200.08 | K | Joback Method |
| vc | 0.380 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 185.94 | J/molxK | 392.48 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 202.04 | J/mol×K | 426.07 | Joback Method |
| cpg | 217.04 | J/mol×K | 459.66 | Joback Method |
| cpg | 230.99 | J/mol×K | 493.26 | Joback Method |
| cpg | 243.98 | J/mol×K | 526.85 | Joback Method |
| cpg | 256.06 | J/mol×K | 560.44 | Joback Method |
| cpg | 267.30 | J/mol×K | 594.03 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=C822935&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |

Legend

| | |
|-----------------|---|
| affp: | Proton affinity |
| basg: | Gas basicity |
| 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 |
| ie: | Ionization energy |
| 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 |

Latest version available from:

<https://www.chemeo.com/cid/27-585-3/Cyclopropane-1-1-ethenyldenebis.pdf>

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