

cis-eudesma-6,11-diene

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| Inchi: | InChI=1S/C15H24/c1-11(2)13-7-9-15(4)8-5-6-12(3)14(15)10-13/h10,12,14H,1,5-9H2,2-4H |
| InchiKey: | GZTVOICLUQHEMR-UHFFFAOYSA-N |
| Formula: | C15H24 |
| SMILES: | <chem>C=C(C)C1=CC2C(C)CCCC2(C)CC1</chem> |
| Mol. weight [g/mol]: | 204.35 |
| CAS: | 194607-93-7 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 234.94 | kJ/mol | Joback Method |
| hf | -75.12 | kJ/mol | Joback Method |
| hfus | 15.49 | kJ/mol | Joback Method |
| hvap | 48.40 | kJ/mol | Joback Method |
| log10ws | -4.87 | | Crippen Method |
| logp | 4.725 | | Crippen Method |
| mcvol | 191.890 | ml/mol | McGowan Method |
| pc | 2032.72 | kPa | Joback Method |
| rinpol | 1484.00 | | NIST Webbook |
| rinpol | 1484.00 | | NIST Webbook |
| rinpol | 1487.00 | | NIST Webbook |
| rinpol | 1492.70 | | NIST Webbook |
| rinpol | 1498.00 | | NIST Webbook |
| ripol | 1692.00 | | NIST Webbook |
| tb | 569.43 | K | Joback Method |
| tc | 794.31 | K | Joback Method |
| tf | 297.83 | K | Joback Method |
| vc | 0.723 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 499.65 | J/molxK | 569.43 | Joback Method |
| cpg | 523.17 | J/molxK | 606.91 | Joback Method |
| cpg | 545.19 | J/molxK | 644.39 | Joback Method |

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|-----|--------|---------|--------|---------------|
| cpg | 565.88 | J/mol×K | 681.87 | Joback Method |
| cpg | 585.40 | J/mol×K | 719.35 | Joback Method |
| cpg | 603.90 | J/mol×K | 756.83 | Joback Method |
| cpg | 621.57 | J/mol×K | 794.31 | Joback Method |

Sources

| | |
|------------------------|---|
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=C194607937&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| 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

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|-----------------|---|
| 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 |
| ripol: | Polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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