

Muurola-4(15),5-diene

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|-----------------------------|---|
| Inchi: | InChI=1S/C15H24/c1-10(2)13-8-6-12(4)14-7-5-11(3)9-15(13)14/h9-10,12-14H,3,5-8H2,1 |
| InchiKey: | RNDFUOKDULDZPR-MELADBBJSA-N |
| Formula: | C15H24 |
| SMILES: | <chem>C=C1C=C2C(C(C)C)CCC(C)C2CC1</chem> |
| Mol. weight [g/mol]: | 204.35 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 211.78 | kJ/mol | Joback Method |
| hf | -127.04 | kJ/mol | Joback Method |
| hfus | 19.70 | kJ/mol | Joback Method |
| hvap | 49.91 | kJ/mol | Joback Method |
| log10ws | -4.63 | | Crippen Method |
| logp | 4.581 | | Crippen Method |
| mcvol | 191.890 | ml/mol | McGowan Method |
| pc | 1918.62 | kPa | Joback Method |
| rinpol | 1467.00 | | NIST Webbook |
| rinpol | 1468.00 | | NIST Webbook |
| tb | 571.35 | K | Joback Method |
| tc | 785.53 | K | Joback Method |
| tf | 288.33 | K | Joback Method |
| vc | 0.721 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 501.64 | J/molxK | 571.35 | Joback Method |
| cpg | 524.63 | J/molxK | 607.05 | Joback Method |
| cpg | 546.31 | J/molxK | 642.74 | Joback Method |
| cpg | 566.71 | J/molxK | 678.44 | Joback Method |
| cpg | 585.88 | J/molxK | 714.14 | Joback Method |
| cpg | 603.88 | J/molxK | 749.83 | Joback Method |
| cpg | 620.74 | J/molxK | 785.53 | Joback Method |
| dvisc | 0.0022043 | Paxs | 288.33 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0013305 | Paxs | 335.50 | Joback Method |
| dvisc | 0.0009095 | Paxs | 382.67 | Joback Method |
| dvisc | 0.0006759 | Paxs | 429.84 | Joback Method |
| dvisc | 0.0005326 | Paxs | 477.01 | Joback Method |
| dvisc | 0.0004381 | Paxs | 524.18 | Joback Method |
| dvisc | 0.0003722 | Paxs | 571.35 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R620898&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

| | |
|----------------------------|---|
| cp_g: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| g_f: | Standard Gibbs free energy of formation |
| h_f: | Enthalpy of formation at standard conditions |
| h_{fus}: | Enthalpy of fusion at standard conditions |
| h_{vap}: | Enthalpy of vaporization at standard conditions |
| log₁₀ws: | Log ₁₀ of Water solubility in mol/l |
| log_p: | Octanol/Water partition coefficient |
| mc_{vol}: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rin_{pol}: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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