

Fokienol

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|-----------------------------|---|
| Other names: | Fokienol [(E,E)-Farnesa-1,6,9,11-tetraen-3-ol] |
| Inchi: | InChI=1S/C17H28O/c1-6-17(5,18)14-10-13-16(4)12-9-7-8-11-15(2)3/h6,8,11,13,18H,1-2 |
| InchiKey: | XCTSDLUOCRRDLZ-YZLWMTBJSA-N |
| Formula: | C17H28O |
| SMILES: | <chem>C=CC(C)(O)CCC=C(C)CCCC=CC(=C)C</chem> |
| Mol. weight [g/mol]: | 248.40 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 277.30 | kJ/mol | Joback Method |
| hf | -89.47 | kJ/mol | Joback Method |
| hfus | 31.68 | kJ/mol | Joback Method |
| hvap | 67.55 | kJ/mol | Joback Method |
| log10ws | -5.73 | | Crippen Method |
| logp | 4.952 | | Crippen Method |
| mcvol | 239.060 | ml/mol | McGowan Method |
| pc | 1548.78 | kPa | Joback Method |
| rinpol | 1576.00 | | NIST Webbook |
| rinpol | 1577.00 | | NIST Webbook |
| rinpol | 1583.00 | | NIST Webbook |
| rinpol | 1583.00 | | NIST Webbook |
| rinpol | 1577.00 | | NIST Webbook |
| rinpol | 1576.00 | | NIST Webbook |
| rinpol | 1568.00 | | NIST Webbook |
| rinpol | 1577.00 | | NIST Webbook |
| ripol | 2170.00 | | NIST Webbook |
| ripol | 2170.00 | | NIST Webbook |
| ripol | 2170.00 | | NIST Webbook |
| ripol | 2177.00 | | NIST Webbook |
| ripol | 2172.00 | | NIST Webbook |
| ripol | 2174.00 | | NIST Webbook |
| ripol | 2187.00 | | NIST Webbook |
| ripol | 2174.00 | | NIST Webbook |
| tb | 678.75 | K | Joback Method |
| tc | 860.47 | K | Joback Method |
| tf | 302.99 | K | Joback Method |
| vc | 0.919 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 659.93 | J/mol×K | 678.75 | Joback Method |
| cpg | 676.03 | J/mol×K | 709.04 | Joback Method |
| cpg | 691.28 | J/mol×K | 739.32 | Joback Method |
| cpg | 705.77 | J/mol×K | 769.61 | Joback Method |
| cpg | 719.55 | J/mol×K | 799.89 | Joback Method |
| cpg | 732.69 | J/mol×K | 830.18 | Joback Method |
| cpg | 745.27 | J/mol×K | 860.47 | Joback Method |

Sources

| | |
|------------------------|---|
| 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=R193801&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307I |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |

Legend

| | |
|-----------------|---|
| 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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<https://www.chemeo.com/cid/79-939-3/Fokienol.pdf>

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