

Phenol, 4-(1,1,2,2,3-pentamethylbutyl)

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|-----------------------------|---|
| Inchi: | InChI=1S/C15H24O/c1-11(2)14(3,4)15(5,6)12-7-9-13(16)10-8-12/h7-11,16H,1-6H3 |
| InchiKey: | UUGGPZQFAKWNT-UHFFFAOYSA-N |
| Formula: | C15H24O |
| SMILES: | CC(C)C(C)(C)C(C)(C)c1ccc(O)cc1 |
| Mol. weight [g/mol]: | 220.35 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | 36.45 | kJ/mol | Joback Method |
| hf | -316.49 | kJ/mol | Joback Method |
| hfus | 16.08 | kJ/mol | Joback Method |
| hvap | 61.29 | kJ/mol | Joback Method |
| log10ws | -3.94 | | Crippen Method |
| logp | 4.352 | | Crippen Method |
| mvol | 204.320 | ml/mol | McGowan Method |
| pc | 2200.01 | kPa | Joback Method |
| rinpol | 1783.00 | | NIST Webbook |
| rinpol | 1783.00 | | NIST Webbook |
| tb | 643.00 | K | Joback Method |
| tc | 872.59 | K | Joback Method |
| tf | 386.79 | K | Joback Method |
| vc | 0.706 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 563.77 | J/molxK | 643.00 | Joback Method |
| cpg | 582.53 | J/molxK | 681.26 | Joback Method |
| cpg | 599.92 | J/molxK | 719.53 | Joback Method |
| cpg | 616.10 | J/molxK | 757.79 | Joback Method |
| cpg | 631.22 | J/molxK | 796.06 | Joback Method |
| cpg | 645.45 | J/molxK | 834.32 | Joback Method |
| cpg | 658.94 | J/molxK | 872.59 | Joback Method |
| dvisc | 0.0019516 | Paxs | 386.79 | Joback Method |

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|-------|-----------|------|--------|---------------|
| dvisc | 0.0005700 | Paxs | 429.49 | Joback Method |
| dvisc | 0.0002080 | Paxs | 472.19 | Joback Method |
| dvisc | 0.0000897 | Paxs | 514.89 | Joback Method |
| dvisc | 0.0000440 | Paxs | 557.60 | Joback Method |
| dvisc | 0.0000239 | Paxs | 600.30 | Joback Method |
| dvisc | 0.0000141 | Paxs | 643.00 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=R592397&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 |
| 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: | Log ₁₀ of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| m_{cvol}: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-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/33-099-6/Phenol-4-1-1-2-2-3-pentamethylbutyl.pdf>

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