

1-Pentanol, 2-amino-4-methyl-, (S)-

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|-----------------------------|---|
| Other names: | (S)-(+)-Leucinol (S)-2-amino-4-methylpentan-1-ol |
| Inchi: | InChI=1S/C6H15NO/c1-5(2)3-6(7)4-8/h5-6,8H,3-4,7H2,1-2H3/t6-/m1/s1 |
| InchiKey: | VPSSPAXIFBTOHY-ZCFIWIBFSA-N |
| Formula: | C6H15NO |
| SMILES: | CC(C)CC(N)CO |
| Mol. weight [g/mol]: | 117.19 |
| CAS: | 7533-40-6 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -75.61 | kJ/mol | Joback Method |
| hf | -296.17 | kJ/mol | Joback Method |
| hfus | 13.53 | kJ/mol | Joback Method |
| hvap | 55.49 | kJ/mol | Joback Method |
| log10ws | -0.90 | | Crippen Method |
| logp | 0.352 | | Crippen Method |
| mvol | 111.250 | ml/mol | McGowan Method |
| pc | 3768.41 | kPa | Joback Method |
| tb | 500.51 | K | Joback Method |
| tc | 680.59 | K | Joback Method |
| tf | 271.46 | K | Joback Method |
| vc | 0.407 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|--------|---------|-----------------|---------------|
| cpg | 261.10 | J/molxK | 500.51 | Joback Method |
| cpg | 271.63 | J/molxK | 530.52 | Joback Method |
| cpg | 281.69 | J/molxK | 560.54 | Joback Method |
| cpg | 291.31 | J/molxK | 590.55 | Joback Method |
| cpg | 300.49 | J/molxK | 620.56 | Joback Method |
| cpg | 309.25 | J/molxK | 650.57 | Joback Method |
| cpg | 317.61 | J/molxK | 680.59 | Joback Method |

Pressure Dependent Properties

| Property code | Value | Unit | Pressure [kPa] | Source |
|---------------|--------|------|----------------|--------------|
| tbrp | 472.20 | K | 102.00 | NIST Webbook |

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=C7533406&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| 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 |
| tb: | Normal Boiling Point Temperature |
| tbrp: | Boiling point at reduced pressure |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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