

D-Leucine

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|----------------------|--|
| Other names: | (R)-(-)-Leucine |
| | (S)-(+)-leucine |
| | (S)-2-amino-4-methylpentanoic acid |
| | (S)-2-amino-4-methylvaleric acid |
| | (S)-leucine |
| | .alpha.-amino-.gamma.-methylvaleric acid |
| | .alpha.-aminoisocaproic acid |
| | 4-methyl-L-norvaline |
| | L-(+)-leucine |
| | L.alpha.-aminoisocaproic acid |
| | L-leucine |
| | Leucine, D- |
| | InChI=1S/C6H13NO2/c1-4(2)3-5(7)6(8)9/h4-5H,3,7H2,1-2H3,(H,8,9)/t5-/m0/s1 |
| Inchi: | |
| InchiKey: | ROHFNLRQFUQHCH-YFKPBYRVSA-N |
| Formula: | C6H13NO2 |
| SMILES: | CC(C)CC(N)C(=O)O |
| Mol. weight [g/mol]: | 131.17 |
| CAS: | 328-38-1 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|-----------------|---------|----------------|
| chs | -3581.40 ± 0.84 | kJ/mol | NIST Webbook |
| gf | -204.53 | kJ/mol | Joback Method |
| hf | -408.75 | kJ/mol | Joback Method |
| hfs | -637.56 ± 0.84 | kJ/mol | NIST Webbook |
| hfus | 15.13 | kJ/mol | Joback Method |
| hvap | 62.24 | kJ/mol | Joback Method |
| log10ws | -0.74 | | Crippen Method |
| logp | 0.444 | | Crippen Method |
| mcvol | 112.820 | ml/mol | McGowan Method |
| pc | 4077.71 | kPa | Joback Method |
| tb | 554.38 | K | Joback Method |
| tc | 743.19 | K | Joback Method |
| tf | 321.39 | K | Joback Method |
| vc | 0.413 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|----------|---------|-----------------|--|
| cpg | 275.71 | J/mol×K | 554.38 | Joback Method |
| cpg | 285.39 | J/mol×K | 585.85 | Joback Method |
| cpg | 294.59 | J/mol×K | 617.32 | Joback Method |
| cpg | 303.34 | J/mol×K | 648.79 | Joback Method |
| cpg | 311.64 | J/mol×K | 680.26 | Joback Method |
| cpg | 319.50 | J/mol×K | 711.72 | Joback Method |
| cpg | 326.94 | J/mol×K | 743.19 | Joback Method |
| psub | 9.00e-04 | kPa | 420.70 | Sublimation and vapour pressure estimation of l-leucine using thermogravimetric analysis |
| psub | 2.90e-03 | kPa | 440.30 | Sublimation and vapour pressure estimation of l-leucine using thermogravimetric analysis |
| psub | 0.04 | kPa | 469.90 | Sublimation and vapour pressure estimation of l-leucine using thermogravimetric analysis |
| psub | 0.53 | kPa | 499.20 | Sublimation and vapour pressure estimation of l-leucine using thermogravimetric analysis |
| psub | 2.16 | kPa | 517.50 | Sublimation and vapour pressure estimation of l-leucine using thermogravimetric analysis |

Sources

Densities, Partial Molar Volumes, and Heat Capacities of Glycine, L-Alanine, and L-Leucine in Aqueous Magnesium Chloride Solutions at Different Temperatures

Solutions at different temperatures in aqueous KCl/KNO₃ systems at different hydrophobic amino acids, peptides, and primary amines with aqueous 3-chloro-1,2-propanediol and 3-chloro-1-propanol: Biophysical studies:

<https://www.doi.org/10.1021/je034168m>

<https://www.doi.org/10.1016/j.jct.2008.07.019>

<https://www.doi.org/10.1016/j.tca.2008.10.023>

<https://www.doi.org/10.1016/j.jct.2010.11.015>

Volumetric behaviour of amino acids and their group contributions in aqueous solution at different temperatures and pressures of different antibiotic drug Chloramphenicol with Electrode and Poly(Vinylidene fluoride) Molal Volumes and Partial Molal (88.15 318.15) Kmol of chloramphenicol in aqueous solution and its activity and group contributions: Thermodynamic considerations of amino acids with sarcosine in presence of Grignard Method Volumetric and

[illegible]

**Volumetric and viscosity properties of
Gripfen in aqueous solutions at T
= (283.15 to 333.15) K:
Intermolecular interactions of
.alpha.-amino acids and glycyl
dipeptides with the drug domiphen
bromide in aqueous solutions analyzed
by volumetric and UV-vis spectroscopy**

chs: Standard solid enthalpy of combustion
cpg: Ideal gas heat capacity
gf: Standard Gibbs free energy of formation

<https://www.doi.org/10.1016/j.ijct.2016.06.018>

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|-----------------|--|
| hf: | Enthalpy of formation at standard conditions |
| hfs: | Solid phase 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 |
| psub: | Sublimation pressure |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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<https://www.cheméo.com/cid/38-802-9/D-Leucine.pdf>

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