

# 3,5-Dimethylheptane, (L)

|                      |                                                                  |
|----------------------|------------------------------------------------------------------|
| Inchi:               | InChI=1S/C9H20/c1-5-8(3)7-9(4)6-2/h8-9H,5-7H2,1-4H3/t8-,9-/m1/s1 |
| InchiKey:            | DZJTZGHZAWTWGA-RKDXNWHRSA-N                                      |
| Formula:             | C9H20                                                            |
| SMILES:              | CCC(C)CC(C)CC                                                    |
| Mol. weight [g/mol]: | 128.26                                                           |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | 20.02   | kJ/mol               | Joback Method  |
| hf            | -239.65 | kJ/mol               | Joback Method  |
| hfus          | 12.02   | kJ/mol               | Joback Method  |
| hvap          | 34.85   | kJ/mol               | Joback Method  |
| log10ws       | -3.11   |                      | Crippen Method |
| logp          | 3.469   |                      | Crippen Method |
| mvol          | 137.670 | ml/mol               | McGowan Method |
| pc            | 2342.82 | kPa                  | Joback Method  |
| rinpol        | 828.60  |                      | NIST Webbook   |
| rinpol        | 834.60  |                      | NIST Webbook   |
| tb            | 404.44  | K                    | Joback Method  |
| tc            | 574.89  | K                    | Joback Method  |
| tf            | 161.19  | K                    | Joback Method  |
| vc            | 0.527   | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 272.52    | J/molxK | 404.44          | Joback Method |
| cpg           | 341.62    | J/molxK | 546.48          | Joback Method |
| cpg           | 328.88    | J/molxK | 518.07          | Joback Method |
| cpg           | 315.61    | J/molxK | 489.67          | Joback Method |
| cpg           | 301.80    | J/molxK | 461.26          | Joback Method |
| cpg           | 287.44    | J/molxK | 432.85          | Joback Method |
| cpg           | 353.86    | J/molxK | 574.89          | Joback Method |
| dvisc         | 0.0002255 | Paxs    | 404.44          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003168 | Paxs | 363.90 | Joback Method |
| dvisc | 0.0004846 | Paxs | 323.36 | Joback Method |
| dvisc | 0.0008373 | Paxs | 282.81 | Joback Method |
| dvisc | 0.0017372 | Paxs | 242.27 | Joback Method |
| dvisc | 0.0048330 | Paxs | 201.73 | Joback Method |
| dvisc | 0.0224972 | Paxs | 161.19 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R300084&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R300084&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>                                 |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                     |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>rinpol:</b>  | Non-polar retention indices                     |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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