

# 4-Heptanol, 3-ethyl-

|                      |                                                          |
|----------------------|----------------------------------------------------------|
| Other names:         | 3-Ethyl-4-heptanol                                       |
| Inchi:               | InChI=1S/C9H20O/c1-4-7-9(10)8(5-2)6-3/h8-10H,4-7H2,1-3H3 |
| InchiKey:            | TVEJPDJKNOLDQ-UHFFFAOYSA-N                               |
| Formula:             | C9H20O                                                   |
| SMILES:              | CCCC(O)C(CC)CC                                           |
| Mol. weight [g/mol]: | 144.25                                                   |
| CAS:                 | 19780-42-8                                               |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -116.80 | kJ/mol  | Joback Method  |
| hf            | -391.88 | kJ/mol  | Joback Method  |
| hfus          | 16.11   | kJ/mol  | Joback Method  |
| hvap          | 51.53   | kJ/mol  | Joback Method  |
| log10ws       | -2.72   |         | Crippen Method |
| logp          | 2.584   |         | Crippen Method |
| mcvol         | 143.540 | ml/mol  | McGowan Method |
| pc            | 2566.29 | kPa     | Joback Method  |
| tb            | 496.62  | K       | Joback Method  |
| tc            | 661.81  | K       | Joback Method  |
| tf            | 222.01  | K       | Joback Method  |
| vc            | 0.546   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 334.77    | J/mol×K | 496.62          | Joback Method |
| cpg           | 394.96    | J/mol×K | 634.28          | Joback Method |
| cpg           | 383.89    | J/mol×K | 606.75          | Joback Method |
| cpg           | 372.35    | J/mol×K | 579.21          | Joback Method |
| cpg           | 360.33    | J/mol×K | 551.68          | Joback Method |
| cpg           | 347.81    | J/mol×K | 524.15          | Joback Method |
| cpg           | 405.56    | J/mol×K | 661.81          | Joback Method |
| dvisc         | 0.0001337 | Pa×s    | 496.62          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0002445 | Pa×s | 450.85 | Joback Method |
| dvisc | 0.0005125 | Pa×s | 405.08 | Joback Method |
| dvisc | 0.0012970 | Pa×s | 359.31 | Joback Method |
| dvisc | 0.0043049 | Pa×s | 313.55 | Joback Method |
| dvisc | 0.0215322 | Pa×s | 267.78 | Joback Method |
| dvisc | 0.2091511 | Pa×s | 222.01 | 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=C19780428&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C19780428&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</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>mccvol:</b>  | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <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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