

# 4-Heptanone, 3-ethyl-

|                      |                                                       |
|----------------------|-------------------------------------------------------|
| Other names:         | 3-Ethyl-4-heptanone                                   |
| Inchi:               | InChI=1S/C9H18O/c1-4-7-9(10)8(5-2)6-3/h8H,4-7H2,1-3H3 |
| InchiKey:            | QMQBUTKELHAKRH-UHFFFAOYSA-N                           |
| Formula:             | C9H18O                                                |
| SMILES:              | CCCC(=O)C(CC)CC                                       |
| Mol. weight [g/mol]: | 142.24                                                |
| CAS:                 | 1528-25-2                                             |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -106.46 | kJ/mol  | Joback Method  |
| hf            | -346.95 | kJ/mol  | Joback Method  |
| hfus          | 17.14   | kJ/mol  | Joback Method  |
| hvap          | 41.99   | kJ/mol  | Joback Method  |
| log10ws       | -2.63   |         | Crippen Method |
| logp          | 2.792   |         | Crippen Method |
| mcvol         | 139.240 | ml/mol  | McGowan Method |
| pc            | 2472.73 | kPa     | Joback Method  |
| tb            | 458.75  | K       | Joback Method  |
| tc            | 636.64  | K       | Joback Method  |
| tf            | 226.12  | K       | Joback Method  |
| vc            | 0.539   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 297.82    | J/mol×K | 458.75          | Joback Method |
| cpg           | 361.70    | J/mol×K | 606.99          | Joback Method |
| cpg           | 350.00    | J/mol×K | 577.34          | Joback Method |
| cpg           | 337.78    | J/mol×K | 547.69          | Joback Method |
| cpg           | 325.02    | J/mol×K | 518.05          | Joback Method |
| cpg           | 311.71    | J/mol×K | 488.40          | Joback Method |
| cpg           | 372.87    | J/mol×K | 636.64          | Joback Method |
| dvisc         | 0.0002699 | Pa×s    | 458.75          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0003608 | Pa×s | 419.98 | Joback Method |
| dvisc | 0.0005116 | Pa×s | 381.21 | Joback Method |
| dvisc | 0.0007850 | Pa×s | 342.44 | Joback Method |
| dvisc | 0.0013439 | Pa×s | 303.66 | Joback Method |
| dvisc | 0.0026929 | Pa×s | 264.89 | Joback Method |
| dvisc | 0.0068479 | Pa×s | 226.12 | 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=C1528252&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C1528252&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>mcvol:</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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