

# 2-propyl hexanal

**Inchi:** InChI=1S/C9H18O/c1-3-5-7-9(8-10)6-4-2/h8-9H,3-7H2,1-2H3  
**InchiKey:** IREORVYCKDQFMD-UHFFFAOYSA-N  
**Formula:** C9H18O  
**SMILES:** CCCCC(C=O)CCC  
**Mol. weight [g/mol]:** 142.24

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -77.06  | kJ/mol  | Joback Method  |
| hf            | -319.95 | kJ/mol  | Joback Method  |
| hfus          | 17.83   | kJ/mol  | Joback Method  |
| hvap          | 41.96   | kJ/mol  | Joback Method  |
| log10ws       | -2.63   |         | Crippen Method |
| logp          | 2.792   |         | Crippen Method |
| mcvol         | 139.240 | ml/mol  | McGowan Method |
| pc            | 2500.00 | kPa     | Joback Method  |
| ripol         | 1279.00 |         | NIST Webbook   |
| ripol         | 1279.00 |         | NIST Webbook   |
| tb            | 453.54  | K       | Joback Method  |
| tc            | 626.89  | K       | Joback Method  |
| tf            | 218.19  | K       | Joback Method  |
| vc            | 0.550   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 298.73    | J/mol×K | 453.54          | Joback Method |
| cpg           | 312.25    | J/mol×K | 482.43          | Joback Method |
| cpg           | 325.23    | J/mol×K | 511.32          | Joback Method |
| cpg           | 337.68    | J/mol×K | 540.22          | Joback Method |
| cpg           | 349.61    | J/mol×K | 569.11          | Joback Method |
| cpg           | 361.03    | J/mol×K | 598.00          | Joback Method |
| cpg           | 371.96    | J/mol×K | 626.89          | Joback Method |
| dvisc         | 0.0079029 | Pa×s    | 218.19          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0029971 | Pa×s | 257.42 | Joback Method |
| dvisc | 0.0014688 | Pa×s | 296.64 | Joback Method |
| dvisc | 0.0008503 | Pa×s | 335.87 | Joback Method |
| dvisc | 0.0005519 | Pa×s | 375.09 | Joback Method |
| dvisc | 0.0003888 | Pa×s | 414.31 | Joback Method |
| dvisc | 0.0002909 | Pa×s | 453.54 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=R328062&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R328062&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>                                     |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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>ripol:</b>   | 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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