

# 6-Nonen-1-ol, (E)-

|                      |                                                                                                      |
|----------------------|------------------------------------------------------------------------------------------------------|
| Other names:         | trans-6-Nonen-1-ol<br>(6E)-6-Nonen-1-ol<br>6-(E)-Nonen-1-ol<br>(E)-6-Nonen-1-ol<br>(E)-non-6-en-1-ol |
| Inchi:               | InChI=1S/C9H18O/c1-2-3-4-5-6-7-8-9-10/h3-4,10H,2,5-9H2,1H3/b4-3+                                     |
| InchiKey:            | XJHRZBIBSSVCEL-ONEGZZNKSA-N                                                                          |
| Formula:             | C9H18O                                                                                               |
| SMILES:              | CCC=CCCCCO                                                                                           |
| Mol. weight [g/mol]: | 142.24                                                                                               |
| CAS:                 | 31502-19-9                                                                                           |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -31.70  | kJ/mol  | Joback Method  |
| hf            | -264.10 | kJ/mol  | Joback Method  |
| hfus          | 23.36   | kJ/mol  | Joback Method  |
| hvap          | 52.27   | kJ/mol  | Joback Method  |
| log10ws       | -2.71   |         | Crippen Method |
| logp          | 2.505   |         | Crippen Method |
| mcvol         | 139.240 | ml/mol  | McGowan Method |
| pc            | 2662.52 | kPa     | Joback Method  |
| ripol         | 1714.00 |         | NIST Webbook   |
| ripol         | 1749.00 |         | NIST Webbook   |
| tb            | 501.66  | K       | Joback Method  |
| tc            | 667.34  | K       | Joback Method  |
| tf            | 246.93  | K       | Joback Method  |
| vc            | 0.538   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 317.11 | J/mol×K | 501.66          | Joback Method |
| cpg           | 329.17 | J/mol×K | 529.27          | Joback Method |

|       |           |         |        |               |
|-------|-----------|---------|--------|---------------|
| cpg   | 340.72    | J/mol×K | 556.89 | Joback Method |
| cpg   | 351.78    | J/mol×K | 584.50 | Joback Method |
| cpg   | 362.36    | J/mol×K | 612.11 | Joback Method |
| cpg   | 372.49    | J/mol×K | 639.73 | Joback Method |
| cpg   | 382.18    | J/mol×K | 667.34 | Joback Method |
| dvisc | 0.0432087 | Pa×s    | 246.93 | Joback Method |
| dvisc | 0.0079294 | Pa×s    | 289.38 | Joback Method |
| dvisc | 0.0022455 | Pa×s    | 331.84 | Joback Method |
| dvisc | 0.0008466 | Pa×s    | 374.30 | Joback Method |
| dvisc | 0.0003894 | Pa×s    | 416.75 | Joback Method |
| dvisc | 0.0002067 | Pa×s    | 459.21 | Joback Method |
| dvisc | 0.0001222 | Pa×s    | 501.66 | Joback Method |

## Sources

|                        |                                                                                                                                               |
|------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| <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>                         |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C31502199&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C31502199&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>                             |

## 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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