

# 2-Nonene, 3-methyl-, (E)-

|                      |                                                                |
|----------------------|----------------------------------------------------------------|
| Inchi:               | InChI=1S/C10H20/c1-4-6-7-8-9-10(3)5-2/h5H,4,6-9H2,1-3H3/b10-5+ |
| InchiKey:            | CMLFMHBAUHOXAV-BJMVGYQFSA-N                                    |
| Formula:             | C10H20                                                         |
| SMILES:              | CC=C(C)CCCCC                                                   |
| Mol. weight [g/mol]: | 140.27                                                         |
| CAS:                 | 17003-99-5                                                     |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 104.99  | kJ/mol  | Joback Method  |
| hf            | -142.30 | kJ/mol  | Joback Method  |
| hfus          | 20.55   | kJ/mol  | Joback Method  |
| hvap          | 37.89   | kJ/mol  | Joback Method  |
| log10ws       | -3.86   |         | Crippen Method |
| logp          | 3.923   |         | Crippen Method |
| mcvol         | 147.460 | ml/mol  | McGowan Method |
| pc            | 2222.89 | kPa     | Joback Method  |
| tb            | 432.24  | K       | Joback Method  |
| tc            | 605.63  | K       | Joback Method  |
| tf            | 183.42  | K       | Joback Method  |
| vc            | 0.577   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 300.23 | J/mol×K | 432.24          | Joback Method |
| cpg           | 315.43 | J/mol×K | 461.14          | Joback Method |
| cpg           | 329.98 | J/mol×K | 490.04          | Joback Method |
| cpg           | 343.91 | J/mol×K | 518.94          | Joback Method |
| cpg           | 357.23 | J/mol×K | 547.83          | Joback Method |
| cpg           | 369.97 | J/mol×K | 576.73          | Joback Method |
| cpg           | 382.16 | J/mol×K | 605.63          | Joback Method |

# Sources

|                        |                                                                                                                                               |
|------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C17003995&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C17003995&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>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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