

# HEPTANOIC ACID, 1,1-DIMETHYLETHYL ESTER

|                      |                                                               |
|----------------------|---------------------------------------------------------------|
| Inchi:               | InChI=1S/C11H22O2/c1-5-6-7-8-9-10(12)13-11(2,3)4/h5-9H2,1-4H3 |
| InchiKey:            | WLUUTGIJNRNHRF-UHFFFAOYSA-N                                   |
| Formula:             | C11H22O2                                                      |
| SMILES:              | CCCCCCC(=O)OC(C)(C)C                                          |
| Mol. weight [g/mol]: | 186.30                                                        |

## Physical Properties

| Property code | Value   | Unit   | Source             |
|---------------|---------|--------|--------------------|
| af            | 0.5782  |        | Cheméo Relay (1.0) |
| hf            | -617.07 | kJ/mol | Cheméo Relay (1.0) |
| hfus          | 19.62   | kJ/mol | Joback Method      |
| hvap          | 56.06   | kJ/mol | Cheméo Relay (1.0) |
| ie            | 9.62    | eV     | Cheméo Relay (1.0) |
| log10ws       | -3.31   |        | Cheméo Relay (1.0) |
| logp          | 3.298   |        | Crippen Method     |
| mcvol         | 173.290 | ml/mol | McGowan Method     |
| pc            | 2049.31 | kPa    | Joback Method      |
| tb            | 459.74  | K      | Cheméo Relay (1.0) |
| tc            | 641.48  | K      | Cheméo Relay (1.0) |
| tf            | 218.04  | K      | Cheméo Relay (1.0) |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 418.15    | J/mol×K | 524.14          | Joback Method |
| cpg           | 434.09    | J/mol×K | 554.22          | Joback Method |
| cpg           | 449.29    | J/mol×K | 584.30          | Joback Method |
| cpg           | 463.79    | J/mol×K | 614.38          | Joback Method |
| cpg           | 477.60    | J/mol×K | 644.47          | Joback Method |
| cpg           | 490.74    | J/mol×K | 674.55          | Joback Method |
| cpg           | 503.23    | J/mol×K | 704.63          | Joback Method |
| dvisc         | 0.0041126 | Pa×s    | 288.31          | Joback Method |
| dvisc         | 0.0018231 | Pa×s    | 327.62          | Joback Method |
| dvisc         | 0.0009620 | Pa×s    | 366.92          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0005745 | Pa×s | 406.23 | Joback Method |
| dvisc | 0.0003757 | Pa×s | 445.53 | Joback Method |
| dvisc | 0.0002633 | Pa×s | 484.84 | Joback Method |
| dvisc | 0.0001946 | Pa×s | 524.14 | Joback Method |

## Sources

|                                  |                                                                                                                                               |
|----------------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>             | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C41084780&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C41084780&amp;Units=SI</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>                         |
| <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>Cheméo Relay (1.0):</b>       |                                                                                                                                               |
| <b>Cheméo Relay (1.0, beta):</b> |                                                                                                                                               |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>af:</b>      | Acentric Factor                                 |
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <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>ie:</b>      | Ionization energy                               |
| <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                   |

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