

# Octanoic acid, 5-methyl, methyl ester

|                             |   |
|-----------------------------|---|
| <b>Other names:</b>         | 4-Methyl-heptanecarboxylicacidmethylester                     |
| <b>Inchi:</b>               | InChI=1S/C10H20O2/c1-4-6-9(2)7-5-8-10(11)12-3/h9H,4-8H2,1-3H3 |
| <b>InchiKey:</b>            | NEOVRVKRDMQDGO-UHFFFAOYSA-N                                   |
| <b>Formula:</b>             | C10H20O2  |
| <b>SMILES:</b>              | CCCC(C)CCCC(=O)OC   |
| <b>Mol. weight [g/mol]:</b> | 172.26  |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -203.04 | kJ/mol  | Joback Method  |
| hf            | -499.81 | kJ/mol  | Joback Method  |
| hfus          | 20.92   | kJ/mol  | Joback Method  |
| hvap          | 46.62   | kJ/mol  | Joback Method  |
| log10ws       | -2.63   |         | Crippen Method |
| logp          | 2.766   |         | Crippen Method |
| mvol          | 159.200 | ml/mol  | McGowan Method |
| pc            | 2216.62 | kPa     | Joback Method  |
| ripol         | 1467.00 |         | NIST Webbook   |
| ripol         | 1467.00 |         | NIST Webbook   |
| tb            | 504.05  | K       | Joback Method  |
| tc            | 679.81  | K       | Joback Method  |
| tf            | 259.62  | K       | Joback Method  |
| vc            | 0.614   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 368.31 | J/molxK | 504.05          | Joback Method |
| cpg           | 382.82 | J/molxK | 533.34          | Joback Method |
| cpg           | 396.77 | J/molxK | 562.64          | Joback Method |
| cpg           | 410.17 | J/molxK | 591.93          | Joback Method |
| cpg           | 423.03 | J/molxK | 621.22          | Joback Method |
| cpg           | 435.35 | J/molxK | 650.52          | Joback Method |
| cpg           | 447.14 | J/molxK | 679.81          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0046674 | Paxs | 259.62 | Joback Method |
| dvisc | 0.0019627 | Paxs | 300.36 | Joback Method |
| dvisc | 0.0010151 | Paxs | 341.10 | Joback Method |
| dvisc | 0.0006043 | Paxs | 381.84 | Joback Method |
| dvisc | 0.0003976 | Paxs | 422.57 | Joback Method |
| dvisc | 0.0002816 | Paxs | 463.31 | Joback Method |
| dvisc | 0.0002109 | Paxs | 504.05 | 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=R2870&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R2870&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>mvol:</b>    | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>riol:</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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