

# 6-Octenoic acid, 3,7-dimethyl-

|                             |   |
|-----------------------------|---|
| <b>Other names:</b>         | Citronellic acid<br>3,7-Dimethyl-6-octenoic acid                            |
| <b>Inchi:</b>               | InChI=1S/C10H18O2/c1-8(2)5-4-6-9(3)7-10(11)12/h5,9H,4,6-7H2,1-3H3,(H,11,12) |
| <b>InchiKey:</b>            | GJWSUKYXUMVMGX-UHFFFAOYSA-N   |
| <b>Formula:</b>             | C10H18O2  |
| <b>SMILES:</b>              | CC(C)=CCCC(C)CC(=O)O  |
| <b>Mol. weight [g/mol]:</b> | 170.25  |
| <b>CAS:</b>                 | 502-47-6  |

## Physical Properties

| Property code | Value   | Unit                 | Source         |
|---------------|---------|----------------------|----------------|
| gf            | -163.19 | kJ/mol               | Joback Method  |
| hf            | -412.39 | kJ/mol               | Joback Method  |
| hfus          | 22.71   | kJ/mol               | Joback Method  |
| hvap          | 60.93   | kJ/mol               | Joback Method  |
| log10ws       | -2.72   |                      | Crippen Method |
| logp          | 2.844   |                      | Crippen Method |
| mcvol         | 154.900 | ml/mol               | McGowan Method |
| pc            | 2608.40 | kPa                  | Joback Method  |
| rinpol        | 1300.00 |                      | NIST Webbook   |
| rinpol        | 1287.00 |                      | NIST Webbook   |
| rinpol        | 1305.00 |                      | NIST Webbook   |
| rinpol        | 1314.00 |                      | NIST Webbook   |
| rinpol        | 1312.00 |                      | NIST Webbook   |
| rinpol        | 1260.00 |                      | NIST Webbook   |
| rinpol        | 1340.80 |                      | NIST Webbook   |
| rinpol        | 1345.40 |                      | NIST Webbook   |
| ripol         | 2255.00 |                      | NIST Webbook   |
| ripol         | 2257.00 |                      | NIST Webbook   |
| ripol         | 2232.00 |                      | NIST Webbook   |
| tb            | 577.85  | K                    | Joback Method  |
| tc            | 756.62  | K                    | Joback Method  |
| tf            | 279.17  | K                    | Joback Method  |
| vc            | 0.596   | m <sup>3</sup> /kmol | Joback Method  |

# Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 383.37 | J/mol×K | 577.85          | Joback Method |
| cpg           | 395.74 | J/mol×K | 607.65          | Joback Method |
| cpg           | 407.51 | J/mol×K | 637.44          | Joback Method |
| cpg           | 418.72 | J/mol×K | 667.24          | Joback Method |
| cpg           | 429.37 | J/mol×K | 697.03          | Joback Method |
| cpg           | 439.51 | J/mol×K | 726.83          | Joback Method |
| cpg           | 449.16 | J/mol×K | 756.62          | Joback Method |
| hvapt         | 68.70  | kJ/mol  | 451.00          | NIST Webbook  |

## Sources

|                        |   |
|------------------------|---|
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</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>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C502476&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C502476&amp;Units=SI</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>hvapt:</b>   | Enthalpy of vaporization at a given temperature |
| <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>rinpol:</b>  | Non-polar retention indices                     |
| <b>ripol:</b>   | Polar retention indices                         |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |

**tf:** Normal melting (fusion) point

**vc:** Critical Volume

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