

# Caprylic anhydride

|                      |                                                                                  |
|----------------------|----------------------------------------------------------------------------------|
| Other names:         | Octanoic acid, anhydride<br>Octanoic anhydride                                   |
| Inchi:               | InChI=1S/C16H30O3/c1-3-5-7-9-11-13-15(17)19-16(18)14-12-10-8-6-4-2/h3-14H2,1-2H3 |
| InchiKey:            | RAFYDKXYXRZODZ-UHFFFAOYSA-N                                                      |
| Formula:             | C16H30O3                                                                         |
| SMILES:              | CCCCCCCC(=O)OC(=O)CCCCCCC                                                        |
| Mol. weight [g/mol]: | 270.41                                                                           |
| CAS:                 | 623-66-5                                                                         |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -279.00 | kJ/mol  | Joback Method  |
| hf            | -730.95 | kJ/mol  | Joback Method  |
| hfus          | 41.58   | kJ/mol  | Joback Method  |
| hvap          | 67.11   | kJ/mol  | Joback Method  |
| log10ws       | -5.16   |         | Crippen Method |
| logp          | 4.777   |         | Crippen Method |
| mcvol         | 245.310 | ml/mol  | McGowan Method |
| pc            | 1429.38 | kPa     | Joback Method  |
| rinpol        | 1878.30 |         | NIST Webbook   |
| tb            | 555.70  | K       | NIST Webbook   |
| tc            | 871.01  | K       | Joback Method  |
| tf            | 392.17  | K       | Joback Method  |
| vc            | 0.962   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 703.61 | J/mol×K | 695.64          | Joback Method |
| cpg           | 720.38 | J/mol×K | 724.87          | Joback Method |
| cpg           | 736.35 | J/mol×K | 754.10          | Joback Method |
| cpg           | 751.54 | J/mol×K | 783.32          | Joback Method |
| cpg           | 765.96 | J/mol×K | 812.55          | Joback Method |
| cpg           | 779.64 | J/mol×K | 841.78          | Joback Method |

|       |           |         |        |               |
|-------|-----------|---------|--------|---------------|
| cpg   | 792.58    | J/mol×K | 871.01 | Joback Method |
| dvisc | 0.0018108 | Pa×s    | 392.17 | Joback Method |
| dvisc | 0.0008918 | Pa×s    | 442.75 | Joback Method |
| dvisc | 0.0005078 | Pa×s    | 493.33 | Joback Method |
| dvisc | 0.0003211 | Pa×s    | 543.90 | Joback Method |
| dvisc | 0.0002195 | Pa×s    | 594.48 | Joback Method |
| dvisc | 0.0001593 | Pa×s    | 645.06 | Joback Method |
| dvisc | 0.0001211 | Pa×s    | 695.64 | Joback Method |

## Pressure Dependent Properties

| Property code | Value  | Unit | Pressure [kPa] | Source       |
|---------------|--------|------|----------------|--------------|
| tbrp          | 459.20 | K    | 2.00           | NIST Webbook |

## Sources

|                 |                                                                                                                                           |
|-----------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| McGowan Method: | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| NIST Webbook:   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C623665&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C623665&amp;Units=SI</a> |
| Crippen Method: | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                 |
| Crippen Method: | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |
| Joback Method:  | <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>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>rinpol:</b>  | Non-polar retention indices                     |

|              |                                   |
|--------------|-----------------------------------|
| <b>tb:</b>   | Normal Boiling Point Temperature  |
| <b>tbrp:</b> | Boiling point at reduced pressure |
| <b>tc:</b>   | Critical Temperature              |
| <b>tf:</b>   | Normal melting (fusion) point     |
| <b>vc:</b>   | Critical Volume                   |

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