

# butyl 7-octenoate

|                      |                                                                      |
|----------------------|----------------------------------------------------------------------|
| Inchi:               | InChI=1S/C12H22O2/c1-3-5-7-8-9-10-12(13)14-11-6-4-2/h3H,1,4-11H2,2H3 |
| InchiKey:            | NFDZCJRATMRYDM-UHFFFAOYSA-N                                          |
| Formula:             | C12H22O2                                                             |
| SMILES:              | C=CCCCCCC(=O)OCCCC                                                   |
| Mol. weight [g/mol]: | 198.30                                                               |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -95.92  | kJ/mol  | Joback Method  |
| hf            | -410.38 | kJ/mol  | Joback Method  |
| hfus          | 28.34   | kJ/mol  | Joback Method  |
| hvap          | 50.79   | kJ/mol  | Joback Method  |
| log10ws       | -3.56   |         | Crippen Method |
| logp          | 3.466   |         | Crippen Method |
| mcvol         | 183.080 | ml/mol  | McGowan Method |
| pc            | 1920.30 | kPa     | Joback Method  |
| ripol         | 1650.00 |         | NIST Webbook   |
| ripol         | 1651.00 |         | NIST Webbook   |
| ripol         | 1650.00 |         | NIST Webbook   |
| ripol         | 1651.00 |         | NIST Webbook   |
| tb            | 546.93  | K       | Joback Method  |
| tc            | 719.58  | K       | Joback Method  |
| tf            | 295.40  | K       | Joback Method  |
| vc            | 0.713   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 444.34 | J/mol×K | 546.93          | Joback Method |
| cpg           | 514.06 | J/mol×K | 690.81          | Joback Method |
| cpg           | 501.31 | J/mol×K | 662.03          | Joback Method |
| cpg           | 487.98 | J/mol×K | 633.26          | Joback Method |
| cpg           | 474.05 | J/mol×K | 604.48          | Joback Method |
| cpg           | 459.51 | J/mol×K | 575.71          | Joback Method |

|       |           |         |        |               |
|-------|-----------|---------|--------|---------------|
| cpg   | 526.24    | J/mol×K | 719.58 | Joback Method |
| dvisc | 0.0001931 | Pa×s    | 546.93 | Joback Method |
| dvisc | 0.0002516 | Pa×s    | 505.01 | Joback Method |
| dvisc | 0.0003439 | Pa×s    | 463.09 | Joback Method |
| dvisc | 0.0005002 | Pa×s    | 421.17 | Joback Method |
| dvisc | 0.0007904 | Pa×s    | 379.24 | Joback Method |
| dvisc | 0.0013996 | Pa×s    | 337.32 | Joback Method |
| dvisc | 0.0029144 | Pa×s    | 295.40 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=R313649&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R313649&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.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</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>ripol:</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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