

# butyl 2-ethylnonanoate

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C15H30O2/c1-4-7-9-10-11-12-14(6-3)15(16)17-13-8-5-2/h14H,4-13H2,1-3H3 |
| <b>InchiKey:</b>            | NBWSUQDYUZUTEG-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C15H30O2   |
| <b>SMILES:</b>              | CCCCCCCC(CC)C(=O)OCCCC   |
| <b>Mol. weight [g/mol]:</b> | 242.40   |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -160.94 | kJ/mol  | Joback Method  |
| hf            | -603.01 | kJ/mol  | Joback Method  |
| hfus          | 33.87   | kJ/mol  | Joback Method  |
| hvap          | 57.75   | kJ/mol  | Joback Method  |
| log10ws       | -4.72   |         | Crippen Method |
| logp          | 4.716   |         | Crippen Method |
| mcvol         | 229.650 | ml/mol  | McGowan Method |
| pc            | 1474.75 | kPa     | Joback Method  |
| ripol         | 1761.00 |         | NIST Webbook   |
| tb            | 618.45  | K       | Joback Method  |
| tc            | 788.43  | K       | Joback Method  |
| tf            | 315.97  | K       | Joback Method  |
| vc            | 0.893   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 621.09    | J/molxK | 618.45          | Joback Method |
| cpg           | 638.80    | J/molxK | 646.78          | Joback Method |
| cpg           | 655.75    | J/molxK | 675.11          | Joback Method |
| cpg           | 671.98    | J/molxK | 703.44          | Joback Method |
| cpg           | 687.48    | J/molxK | 731.77          | Joback Method |
| cpg           | 702.28    | J/molxK | 760.10          | Joback Method |
| cpg           | 716.38    | J/molxK | 788.43          | Joback Method |
| dvisc         | 0.0034492 | Paxs    | 315.97          | Joback Method |
| dvisc         | 0.0013649 | Paxs    | 366.38          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0006759 | Paxs | 416.80 | Joback Method |
| dvisc | 0.0003895 | Paxs | 467.21 | Joback Method |
| dvisc | 0.0002499 | Paxs | 517.62 | Joback Method |
| dvisc | 0.0001735 | Paxs | 568.04 | Joback Method |
| dvisc | 0.0001278 | Paxs | 618.45 | 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=R493275&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=R493275&amp;Units=SI</a> |
| <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>                         |

## 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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