

# Adipic acid, heptyl 3-methylbutyl ester

|                      |                                                                                  |
|----------------------|----------------------------------------------------------------------------------|
| Inchi:               | InChI=1S/C18H34O4/c1-4-5-6-7-10-14-21-17(19)11-8-9-12-18(20)22-15-13-16(2)3/h16H |
| InchiKey:            | RUWXCFUVXBMOFT-UHFFFAOYSA-N                                                      |
| Formula:             | C18H34O4                                                                         |
| SMILES:              | CCCCCCCCOC(=O)CCCCC(=O)OCCC(C)C                                                  |
| Mol. weight [g/mol]: | 314.46                                                                           |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -369.60 | kJ/mol  | Joback Method  |
| hf            | -909.73 | kJ/mol  | Joback Method  |
| hfus          | 44.43   | kJ/mol  | Joback Method  |
| hvap          | 73.59   | kJ/mol  | Joback Method  |
| log10ws       | -4.84   |         | Crippen Method |
| logp          | 4.650   |         | Crippen Method |
| mcvol         | 279.360 | ml/mol  | McGowan Method |
| pc            | 1234.61 | kPa     | Joback Method  |
| rinpol        | 2108.00 |         | NIST Webbook   |
| rinpol        | 2108.00 |         | NIST Webbook   |
| tb            | 763.38  | K       | Joback Method  |
| tc            | 943.41  | K       | Joback Method  |
| tf            | 421.94  | K       | Joback Method  |
| vc            | 1.085   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 849.98    | J/mol×K | 763.38          | Joback Method |
| cpg           | 867.53    | J/mol×K | 793.39          | Joback Method |
| cpg           | 884.14    | J/mol×K | 823.39          | Joback Method |
| cpg           | 899.82    | J/mol×K | 853.40          | Joback Method |
| cpg           | 914.60    | J/mol×K | 883.40          | Joback Method |
| cpg           | 928.46    | J/mol×K | 913.41          | Joback Method |
| cpg           | 941.44    | J/mol×K | 943.41          | Joback Method |
| dvisc         | 0.0012393 | Pa×s    | 421.94          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0005664 | Pa×s | 478.85 | Joback Method |
| dvisc | 0.0003058 | Pa×s | 535.75 | Joback Method |
| dvisc | 0.0001858 | Pa×s | 592.66 | Joback Method |
| dvisc | 0.0001232 | Pa×s | 649.57 | Joback Method |
| dvisc | 0.0000873 | Pa×s | 706.47 | Joback Method |
| dvisc | 0.0000651 | Pa×s | 763.38 | 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=U348969&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U348969&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>                         |

## 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>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

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