

# Malonic acid, 3-methylpentyl octyl ester

|                      |                                                                                   |
|----------------------|-----------------------------------------------------------------------------------|
| Inchi:               | InChI=1S/C17H32O4/c1-4-6-7-8-9-10-12-20-16(18)14-17(19)21-13-11-15(3)5-2/h15H,4-1 |
| InchiKey:            | XWOGFISJHITWKS-UHFFFAOYSA-N                                                       |
| Formula:             | C17H32O4                                                                          |
| SMILES:              | CCCCCCCCOC(=O)CC(=O)OCCC(C)CC                                                     |
| Mol. weight [g/mol]: | 300.43                                                                            |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -378.02 | kJ/mol  | Joback Method  |
| hf            | -889.09 | kJ/mol  | Joback Method  |
| hfus          | 41.84   | kJ/mol  | Joback Method  |
| hvap          | 71.36   | kJ/mol  | Joback Method  |
| log10ws       | -4.42   |         | Crippen Method |
| logp          | 4.260   |         | Crippen Method |
| mcvol         | 265.270 | ml/mol  | McGowan Method |
| pc            | 1322.31 | kPa     | Joback Method  |
| rinpol        | 1986.00 |         | NIST Webbook   |
| rinpol        | 1986.00 |         | NIST Webbook   |
| tb            | 740.50  | K       | Joback Method  |
| tc            | 919.30  | K       | Joback Method  |
| tf            | 410.67  | K       | Joback Method  |
| vc            | 1.030   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 791.52    | J/mol×K | 740.50          | Joback Method |
| cpg           | 808.68    | J/mol×K | 770.30          | Joback Method |
| cpg           | 824.95    | J/mol×K | 800.10          | Joback Method |
| cpg           | 840.35    | J/mol×K | 829.90          | Joback Method |
| cpg           | 854.88    | J/mol×K | 859.70          | Joback Method |
| cpg           | 868.56    | J/mol×K | 889.50          | Joback Method |
| cpg           | 881.39    | J/mol×K | 919.30          | Joback Method |
| dvisc         | 0.0013763 | Pa×s    | 410.67          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0006360 | Pa×s | 465.64 | Joback Method |
| dvisc | 0.0003459 | Pa×s | 520.61 | Joback Method |
| dvisc | 0.0002113 | Pa×s | 575.59 | Joback Method |
| dvisc | 0.0001407 | Pa×s | 630.56 | Joback Method |
| dvisc | 0.0001000 | Pa×s | 685.53 | Joback Method |
| dvisc | 0.0000748 | Pa×s | 740.50 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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>                         |
| <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=U348282&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U348282&amp;Units=SI</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>mccvol:</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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