

# Sebacic acid, dodecyl heptyl ester

|                             |  |
|-----------------------------|--|
| <b>Inchi:</b>               | InChI=1S/C29H56O4/c1-3-5-7-9-10-11-12-15-19-23-27-33-29(31)25-21-17-14-13-16-20- |
| <b>InchiKey:</b>            | ZZNQTMOQTNOYES-UHFFFAOYSA-N  |
| <b>Formula:</b>             | C29H56O4   |
| <b>SMILES:</b>              | CCCCCCCCCCCCOC(=O)CCCCCCCCC(=O)OCCCCCCC  |
| <b>Mol. weight [g/mol]:</b> | 468.75   |

## Physical Properties

| Property code | Value    | Unit                 | Source         |
|---------------|----------|----------------------|----------------|
| gf            | -274.54  | kJ/mol               | Joback Method  |
| hf            | -1131.49 | kJ/mol               | Joback Method  |
| hfus          | 76.44    | kJ/mol               | Joback Method  |
| hvap          | 98.46    | kJ/mol               | Joback Method  |
| log10ws       | -9.69    |                      | Crippen Method |
| logp          | 9.085    |                      | Crippen Method |
| mvol          | 434.350  | ml/mol               | McGowan Method |
| pc            | 654.10   | kPa                  | Joback Method  |
| rinpol        | 3334.00  |                      | NIST Webbook   |
| rinpol        | 3334.00  |                      | NIST Webbook   |
| tb            | 1015.50  | K                    | Joback Method  |
| tc            | 1268.62  | K                    | Joback Method  |
| tf            | 560.91   | K                    | Joback Method  |
| vc            | 1.708    | m <sup>3</sup> /kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1539.74   | J/molxK | 1015.50         | Joback Method |
| cpg           | 1563.27   | J/molxK | 1057.69         | Joback Method |
| cpg           | 1584.52   | J/molxK | 1099.87         | Joback Method |
| cpg           | 1603.59   | J/molxK | 1142.06         | Joback Method |
| cpg           | 1620.57   | J/molxK | 1184.25         | Joback Method |
| cpg           | 1635.54   | J/molxK | 1226.44         | Joback Method |
| cpg           | 1648.61   | J/molxK | 1268.62         | Joback Method |
| dvisc         | 0.0002741 | Paxs    | 560.91          | Joback Method |

|       |           |      |         |               |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0001230 | Paxs | 636.67  | Joback Method |
| dvisc | 0.0000654 | Paxs | 712.44  | Joback Method |
| dvisc | 0.0000393 | Paxs | 788.20  | Joback Method |
| dvisc | 0.0000258 | Paxs | 863.97  | Joback Method |
| dvisc | 0.0000181 | Paxs | 939.73  | Joback Method |
| dvisc | 0.0000134 | Paxs | 1015.50 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <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>                         |
| <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=U354309&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U354309&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>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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