

# Adipic acid, 2-decyl tetradecyl ester

**Inchi:** InChI=1S/C30H58O4/c1-4-6-8-10-12-13-14-15-16-17-19-23-27-33-29(31)25-21-22-26-30  
**InchiKey:** JKKPWPAWRLYBIN-UHFFFAOYSA-N  
**Formula:** C30H58O4  
**SMILES:** CCCCCCCCCCCCCCOC(=O)CCCCC(=O)OC(C)CCCCCCCC  
**Mol. weight [g/mol]:** 482.78

## Physical Properties

| Property code | Value    | Unit    | Source         |
|---------------|----------|---------|----------------|
| gf            | -268.56  | kJ/mol  | Joback Method  |
| hf            | -1157.41 | kJ/mol  | Joback Method  |
| hfus          | 75.51    | kJ/mol  | Joback Method  |
| hvap          | 100.30   | kJ/mol  | Joback Method  |
| log10ws       | -10.22   |         | Crippen Method |
| logp          | 9.473    |         | Crippen Method |
| mcvol         | 448.440  | ml/mol  | McGowan Method |
| pc            | 625.63   | kPa     | Joback Method  |
| rinpol        | 3218.00  |         | NIST Webbook   |
| tb            | 1037.94  | K       | Joback Method  |
| tc            | 1301.70  | K       | Joback Method  |
| tf            | 557.18   | K       | Joback Method  |
| vc            | 1.758    | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1605.40   | J/mol×K | 1037.94         | Joback Method |
| cpg           | 1629.35   | J/mol×K | 1081.90         | Joback Method |
| cpg           | 1650.81   | J/mol×K | 1125.86         | Joback Method |
| cpg           | 1669.88   | J/mol×K | 1169.82         | Joback Method |
| cpg           | 1686.68   | J/mol×K | 1213.78         | Joback Method |
| cpg           | 1701.32   | J/mol×K | 1257.74         | Joback Method |
| cpg           | 1713.90   | J/mol×K | 1301.70         | Joback Method |
| dvisc         | 0.0002682 | Pa×s    | 557.18          | Joback Method |
| dvisc         | 0.0001110 | Pa×s    | 637.31          | Joback Method |

|       |           |      |         |               |
|-------|-----------|------|---------|---------------|
| dvisc | 0.0000560 | Pa×s | 717.43  | Joback Method |
| dvisc | 0.0000324 | Pa×s | 797.56  | Joback Method |
| dvisc | 0.0000207 | Pa×s | 877.69  | Joback Method |
| dvisc | 0.0000143 | Pa×s | 957.81  | Joback Method |
| dvisc | 0.0000104 | Pa×s | 1037.94 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=U353607&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U353607&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>                                 |

## 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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