

# Sebacic acid, dodecyl trans-3-hexenyl ester

**Inchi:** InChI=1S/C28H52O4/c1-3-5-7-9-10-11-12-15-18-22-26-32-28(30)24-20-17-14-13-16-19-  
**InchiKey:** PZKBBINKHOLBPZ-SOFGYWHQSA-N  
**Formula:** C28H52O4  
**SMILES:** CCC=CCCOC(=O)CCCCCCCCC(=O)OCCCCCCCCCCCCC  
**Mol. weight [g/mol]:** 452.71

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -202.74 | kJ/mol  | Joback Method  |
| hf            | -993.63 | kJ/mol  | Joback Method  |
| hfus          | 74.05   | kJ/mol  | Joback Method  |
| hvap          | 96.19   | kJ/mol  | Joback Method  |
| log10ws       | -9.12   |         | Crippen Method |
| logp          | 8.471   |         | Crippen Method |
| mcvol         | 415.960 | ml/mol  | McGowan Method |
| pc            | 706.58  | kPa     | Joback Method  |
| rinpol        | 3173.00 |         | NIST Webbook   |
| tb            | 996.78  | K       | Joback Method  |
| tc            | 1234.53 | K       | Joback Method  |
| tf            | 544.56  | K       | Joback Method  |
| vc            | 1.631   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1446.46   | J/mol×K | 996.78          | Joback Method |
| cpg           | 1540.23   | J/mol×K | 1194.90         | Joback Method |
| cpg           | 1524.71   | J/mol×K | 1155.28         | Joback Method |
| cpg           | 1507.66   | J/mol×K | 1115.65         | Joback Method |
| cpg           | 1489.00   | J/mol×K | 1076.03         | Joback Method |
| cpg           | 1468.63   | J/mol×K | 1036.40         | Joback Method |
| cpg           | 1554.32   | J/mol×K | 1234.53         | Joback Method |
| dvisc         | 0.0000139 | Pa×s    | 996.78          | Joback Method |
| dvisc         | 0.0000187 | Pa×s    | 921.41          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000266 | Pa×s | 846.04 | Joback Method |
| dvisc | 0.0000407 | Pa×s | 770.67 | Joback Method |
| dvisc | 0.0000681 | Pa×s | 695.30 | Joback Method |
| dvisc | 0.0001291 | Pa×s | 619.93 | Joback Method |
| dvisc | 0.0002922 | Pa×s | 544.56 | 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=U356068&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U356068&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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