

# Succinic acid, 6-methylhept-2-yl octadecyl ester

**Inchi:** InChI=1S/C30H58O4/c1-5-6-7-8-9-10-11-12-13-14-15-16-17-18-19-20-26-33-29(31)24-25  
**InchiKey:** UXZZDUCREHJGTM-UHFFFAOYSA-N  
**Formula:** C30H58O4  
**SMILES:** CCCCCCCCCCCCCCCCCCOC(=O)CCC(=O)OC(C)CCCC(C)C  
**Mol. weight [g/mol]:** 482.78

## Physical Properties

| Property code | Value    | Unit    | Source         |
|---------------|----------|---------|----------------|
| gf            | -271.00  | kJ/mol  | Joback Method  |
| hf            | -1162.69 | kJ/mol  | Joback Method  |
| hfus          | 71.98    | kJ/mol  | Joback Method  |
| hvap          | 99.91    | kJ/mol  | Joback Method  |
| log10ws       | -9.98    |         | Crippen Method |
| logp          | 9.329    |         | Crippen Method |
| mcvol         | 448.440  | ml/mol  | McGowan Method |
| pc            | 628.14   | kPa     | Joback Method  |
| rinpol        | 3205.00  |         | NIST Webbook   |
| rinpol        | 3205.00  |         | NIST Webbook   |
| tb            | 1037.50  | K       | Joback Method  |
| tc            | 1297.92  | K       | Joback Method  |
| tf            | 542.18   | K       | Joback Method  |
| vc            | 1.752    | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1605.60   | J/molxK | 1037.50         | Joback Method |
| cpg           | 1700.04   | J/molxK | 1254.52         | Joback Method |
| cpg           | 1685.63   | J/molxK | 1211.11         | Joback Method |
| cpg           | 1669.09   | J/molxK | 1167.71         | Joback Method |
| cpg           | 1650.31   | J/molxK | 1124.31         | Joback Method |
| cpg           | 1629.18   | J/molxK | 1080.90         | Joback Method |
| cpg           | 1712.42   | J/molxK | 1297.92         | Joback Method |
| dvisc         | 0.0000094 | Paxs    | 1037.50         | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000131 | Paxs | 954.95 | Joback Method |
| dvisc | 0.0000194 | Paxs | 872.39 | Joback Method |
| dvisc | 0.0000312 | Paxs | 789.84 | Joback Method |
| dvisc | 0.0000558 | Paxs | 707.29 | Joback Method |
| dvisc | 0.0001167 | Paxs | 624.73 | Joback Method |
| dvisc | 0.0003054 | Paxs | 542.18 | Joback Method |

## Sources

|                        |   |
|------------------------|---|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U381377&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U381377&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>                         |
| <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>                     |

## 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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