

# Succinic acid, hept-2-yl 3-heptyl ester

**Inchi:** InChI=1S/C18H34O4/c1-5-8-10-11-15(4)21-17(19)13-14-18(20)22-16(7-3)12-9-6-2/h15-17  
**InchiKey:** PXAXWSSJNOAODR-UHFFFAOYSA-N  
**Formula:** C18H34O4  
**SMILES:** CCCCCC(C)OC(=O)CCC(=O)OC(CC)CCCC  
**Mol. weight [g/mol]:** 314.46

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -372.04 | kJ/mol  | Joback Method  |
| hf            | -915.01 | kJ/mol  | Joback Method  |
| hfus          | 40.90   | kJ/mol  | Joback Method  |
| hvap          | 73.20   | kJ/mol  | Joback Method  |
| log10ws       | -5.31   |         | Crippen Method |
| logp          | 4.791   |         | Crippen Method |
| mcvol         | 279.360 | ml/mol  | McGowan Method |
| pc            | 1241.58 | kPa     | Joback Method  |
| rinpol        | 1933.00 |         | NIST Webbook   |
| rinpol        | 1933.00 |         | NIST Webbook   |
| tb            | 762.94  | K       | Joback Method  |
| tc            | 944.42  | K       | Joback Method  |
| tf            | 406.94  | K       | Joback Method  |
| vc            | 1.079   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 850.46    | J/mol×K | 762.94          | Joback Method |
| cpg           | 868.17    | J/mol×K | 793.19          | Joback Method |
| cpg           | 884.91    | J/mol×K | 823.43          | Joback Method |
| cpg           | 900.70    | J/mol×K | 853.68          | Joback Method |
| cpg           | 915.55    | J/mol×K | 883.93          | Joback Method |
| cpg           | 929.48    | J/mol×K | 914.17          | Joback Method |
| cpg           | 942.50    | J/mol×K | 944.42          | Joback Method |
| dvisc         | 0.0015163 | Pa×s    | 406.94          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0006283 | Pa×s | 466.27 | Joback Method |
| dvisc | 0.0003176 | Pa×s | 525.61 | Joback Method |
| dvisc | 0.0001844 | Pa×s | 584.94 | Joback Method |
| dvisc | 0.0001183 | Pa×s | 644.27 | Joback Method |
| dvisc | 0.0000818 | Pa×s | 703.61 | Joback Method |
| dvisc | 0.0000599 | Pa×s | 762.94 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U390600&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U390600&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>cp<sub>g</sub>:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>              | Dynamic viscosity                               |
| <b>g<sub>f</sub>:</b>      | Standard Gibbs free energy of formation         |
| <b>h<sub>f</sub>:</b>      | Enthalpy of formation at standard conditions    |
| <b>h<sub>fus</sub>:</b>    | Enthalpy of fusion at standard conditions       |
| <b>h<sub>vap</sub>:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log<sub>10ws</sub>:</b> | Log10 of Water solubility in mol/l              |
| <b>log<sub>p</sub>:</b>    | Octanol/Water partition coefficient             |
| <b>mc<sub>vol</sub>:</b>   | McGowan's characteristic volume                 |
| <b>p<sub>c</sub>:</b>      | Critical Pressure                               |
| <b>rin<sub>pol</sub>:</b>  | Non-polar retention indices                     |
| <b>t<sub>b</sub>:</b>      | Normal Boiling Point Temperature                |
| <b>t<sub>c</sub>:</b>      | Critical Temperature                            |
| <b>t<sub>f</sub>:</b>      | Normal melting (fusion) point                   |
| <b>v<sub>c</sub>:</b>      | Critical Volume                                 |

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