

# Glutaric acid, 6-ethyloct-3-yl isoheptyl ester

**Inchi:** InChI=1S/C21H40O4/c1-6-18(7-2)14-15-19(8-3)25-21(23)13-9-12-20(22)24-16-10-11-17  
**InchiKey:** OCLXJBIAOOIBMZ-UHFFFAOYSA-N  
**Formula:** C21H40O4  
**SMILES:** CCC(CC)CCC(CC)OC(=O)CCCC(=O)OCCCC(C)C  
**Mol. weight [g/mol]:** 356.54

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | -349.22 | kJ/mol  | Joback Method  |
| hf            | -982.21 | kJ/mol  | Joback Method  |
| hfus          | 45.15   | kJ/mol  | Joback Method  |
| hvap          | 79.49   | kJ/mol  | Joback Method  |
| log10ws       | -5.97   |         | Crippen Method |
| logp          | 5.674   |         | Crippen Method |
| mcvol         | 321.630 | ml/mol  | McGowan Method |
| pc            | 1028.60 | kPa     | Joback Method  |
| rinpol        | 2278.00 |         | NIST Webbook   |
| tb            | 831.14  | K       | Joback Method  |
| tc            | 1020.11 | K       | Joback Method  |
| tf            | 425.75  | K       | Joback Method  |
| vc            | 1.242   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 1031.67   | J/mol×K | 831.14          | Joback Method |
| cpg           | 1114.48   | J/mol×K | 988.62          | Joback Method |
| cpg           | 1100.16   | J/mol×K | 957.12          | Joback Method |
| cpg           | 1084.74   | J/mol×K | 925.63          | Joback Method |
| cpg           | 1068.19   | J/mol×K | 894.13          | Joback Method |
| cpg           | 1050.51   | J/mol×K | 862.64          | Joback Method |
| cpg           | 1127.72   | J/mol×K | 1020.11         | Joback Method |
| dvisc         | 0.0000355 | Pa×s    | 831.14          | Joback Method |
| dvisc         | 0.0000496 | Pa×s    | 763.58          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0000739 | Pa×s | 696.01 | Joback Method |
| dvisc | 0.0001201 | Pa×s | 628.44 | Joback Method |
| dvisc | 0.0002191 | Pa×s | 560.88 | Joback Method |
| dvisc | 0.0004714 | Pa×s | 493.32 | Joback Method |
| dvisc | 0.0012935 | Pa×s | 425.75 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <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=U358219&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U358219&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>                                     |

## 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                                 |

Latest version available from:

<https://www.chemeo.com/cid/23-620-7/Glutaric-acid-6-ethyloct-3-yl-isohexyl-ester.pdf>

Generated by Cheméo on 2025-12-05 11:04:51.508831356 +0000 UTC m=+4680889.038872013.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.