

Glutaric acid, 3-methylbut-2-yl isopropyl ester

Inchi:	InChI=1S/C13H24O4/c1-9(2)11(5)17-13(15)8-6-7-12(14)16-10(3)4/h9-11H,6-8H2,1-5H3
InchiKey:	WITDUQPWOBYWRC-UHFFFAOYSA-N
Formula:	C13H24O4
SMILES:	CC(C)OC(=O)CCCC(=O)OC(C)C(C)C
Mol. weight [g/mol]:	244.33

Physical Properties

Property code	Value	Unit	Source
gf	-416.58	kJ/mol	Joback Method
hf	-817.09	kJ/mol	Joback Method
hfus	24.43	kJ/mol	Joback Method
hvap	61.68	kJ/mol	Joback Method
log10ws	-2.97		Crippen Method
logp	2.696		Crippen Method
mcvol	208.910	ml/mol	McGowan Method
pc	1810.77	kPa	Joback Method
rinpol	1478.00		NIST Webbook
rinpol	1478.00		NIST Webbook
tb	648.10	K	Joback Method
tc	832.83	K	Joback Method
tf	335.59	K	Joback Method
vc	0.793	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	569.70	J/mol×K	648.10	Joback Method
cpg	585.64	J/mol×K	678.89	Joback Method
cpg	600.82	J/mol×K	709.68	Joback Method
cpg	615.23	J/mol×K	740.46	Joback Method
cpg	628.87	J/mol×K	771.25	Joback Method
cpg	641.76	J/mol×K	802.04	Joback Method
cpg	653.89	J/mol×K	832.83	Joback Method
dvisc	0.0033428	Pa×s	335.59	Joback Method

dvisc	0.0012868	Pa×s	387.68	Joback Method
dvisc	0.0006210	Pa×s	439.76	Joback Method
dvisc	0.0003497	Pa×s	491.85	Joback Method
dvisc	0.0002198	Pa×s	543.93	Joback Method
dvisc	0.0001499	Pa×s	596.02	Joback Method
dvisc	0.0001087	Pa×s	648.10	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U393369&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/90-836-4/Glutaric-acid-3-methylbut-2-yl-isopropyl-ester.pdf>

Generated by Cheméo on 2025-12-06 02:40:01.192065035 +0000 UTC m=+4736998.722105689.

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