

Glutaric acid, dodecyl 4-nitrophenyl ester

Inchi:	InChI=1S/C23H35NO6/c1-2-3-4-5-6-7-8-9-10-11-19-29-22(25)13-12-14-23(26)30-21-17-
InchiKey:	USZRSKJZSDAWIC-UHFFFAOYSA-N
Formula:	C23H35NO6
SMILES:	CCCCCCCCCCCCOC(=O)CCCC(=O)Oc1ccc([N+](=O)[O-])cc1
Mol. weight [g/mol]:	421.53

Physical Properties

Property code	Value	Unit	Source
gf	-186.73	kJ/mol	Joback Method
hf	-793.35	kJ/mol	Joback Method
hfus	65.91	kJ/mol	Joback Method
hvap	104.63	kJ/mol	Joback Method
log10ws	-7.58		Crippen Method
logp	6.135		Crippen Method
mcvol	343.470	ml/mol	McGowan Method
pc	1106.68	kPa	Joback Method
rinpola	3341.00		NIST Webbook
tb	1061.72	K	Joback Method
tc	1300.03	K	Joback Method
tf	675.84	K	Joback Method
vc	1.345	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1169.01	J/mol×K	1061.72	Joback Method
cpg	1181.96	J/mol×K	1101.44	Joback Method
cpg	1193.35	J/mol×K	1141.16	Joback Method
cpg	1203.22	J/mol×K	1180.88	Joback Method
cpg	1211.64	J/mol×K	1220.59	Joback Method
cpg	1218.67	J/mol×K	1260.31	Joback Method
cpg	1224.35	J/mol×K	1300.03	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U359205&Units=SI

Legend

cpg:	Ideal gas heat capacity
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
mccvol:	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/51-377-7/Glutaric-acid-dodecyl-4-nitrophenyl-ester.pdf>

Generated by Cheméo on 2024-04-18 04:44:28.175102822 +0000 UTC m=+15704717.095680138.

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