

Glutaric acid, dodec-2-en-1-yl 4-chloro-2-methoxyphenyl ester

Inchi:	InChI=1S/C24H35ClO5/c1-3-4-5-6-7-8-9-10-11-12-18-29-23(26)14-13-15-24(27)30-21-17
InchiKey:	HLZDGIOMLLNGET-VAWYXSNFSA-N
Formula:	C24H35ClO5
SMILES:	CCCCCCCCC=CCOC(=O)CCCC(=O)Oc1ccc(Cl)cc1OC
Mol. weight [g/mol]:	438.99

Physical Properties

Property code	Value	Unit	Source
gf	-260.20	kJ/mol	Joback Method
hf	-845.44	kJ/mol	Joback Method
hfus	62.34	kJ/mol	Joback Method
hvap	97.68	kJ/mol	Joback Method
log10ws	-7.59		Crippen Method
logp	6.664		Crippen Method
mvol	353.950	ml/mol	McGowan Method
pc	1020.08	kPa	Joback Method
rinpol	3161.00		NIST Webbook
rinpol	3161.00		NIST Webbook
tb	1001.75	K	Joback Method
tc	1226.43	K	Joback Method
tf	603.09	K	Joback Method
vc	1.367	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1144.85	J/molxK	1001.75	Joback Method
cpg	1203.85	J/molxK	1188.99	Joback Method
cpg	1194.85	J/molxK	1151.54	Joback Method
cpg	1184.49	J/molxK	1114.09	Joback Method
cpg	1172.72	J/molxK	1076.64	Joback Method
cpg	1159.53	J/molxK	1039.20	Joback Method
cpg	1211.54	J/molxK	1226.43	Joback Method
dvisc	0.0000172	Paxs	1001.75	Joback Method

dvisc	0.0000221	Paxs	935.31	Joback Method
dvisc	0.0000295	Paxs	868.86	Joback Method
dvisc	0.0000411	Paxs	802.42	Joback Method
dvisc	0.0000610	Paxs	735.98	Joback Method
dvisc	0.0000980	Paxs	669.53	Joback Method
dvisc	0.0001745	Paxs	603.09	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U393915&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
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/86-763-0/Glutaric-acid-dodec-2-en-1-yl-4-chloro-2-methoxyphenyl-ester.pdf>

Generated by Cheméo on 2024-04-29 14:23:52.548465811 +0000 UTC m=+16689881.469043126.

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