

GA24 methyl ester

Inchi: InChI=1S/C22H30O5/c1-13-10-22-11-14(13)6-7-15(22)21(12-23)9-5-8-20(2,19(25)27-4)1
InchiKey: CDZKHHKZGGMPOC-NJXVUKDQSA-N
Formula: C22H30O5
SMILES: C=C1CC23CC1CCC2C1(C=O)CCCC(C)(C(=O)OC)C1C3C(=O)OC
Mol. weight [g/mol]: 374.47

Physical Properties

Property code	Value	Unit	Source
gf	-212.82	kJ/mol	Joback Method
hf	-730.93	kJ/mol	Joback Method
hfus	30.00	kJ/mol	Joback Method
hvap	85.55	kJ/mol	Joback Method
log10ws	-3.78		Crippen Method
logp	3.317		Crippen Method
mcvol	289.550	ml/mol	McGowan Method
pc	1582.23	kPa	Joback Method
rinpol	2475.00		NIST Webbook
rinpol	2442.00		NIST Webbook
rinpol	2442.00		NIST Webbook
rinpol	2465.00		NIST Webbook
rinpol	2465.00		NIST Webbook
rinpol	2475.00		NIST Webbook
rinpol	2446.00		NIST Webbook
tb	929.64	K	Joback Method
tc	1164.39	K	Joback Method
tf	657.88	K	Joback Method
vc	1.111	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1041.61	J/mol×K	929.64	Joback Method
cpg	1070.38	J/mol×K	968.77	Joback Method
cpg	1100.77	J/mol×K	1007.89	Joback Method

cpg	1133.19	J/mol×K	1047.02	Joback Method
cpg	1168.10	J/mol×K	1086.14	Joback Method
cpg	1205.91	J/mol×K	1125.27	Joback Method
cpg	1247.07	J/mol×K	1164.39	Joback Method

Sources

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=R493433&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071

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
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpola:	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/66-916-2/GA24-methyl-ester.pdf>

Generated by Cheméo on 2022-09-24 23:44:50.202491373 +0000 UTC m=+103028.382288423.

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