

Adipic acid, dodecyl pent-4-en-2-yl ester

Inchi:	InChI=1S/C23H42O4/c1-4-6-7-8-9-10-11-12-13-16-20-26-22(24)18-14-15-19-23(25)27-2
InchiKey:	UUUOTHBLBKQGPA-UHFFFAOYSA-N
Formula:	C23H42O4
SMILES:	C=CCC(C)OC(=O)CCCC(=O)OCCCCCCCCCCCC
Mol. weight [g/mol]:	382.58

Physical Properties

Property code	Value	Unit	Source
gf	-239.66	kJ/mol	Joback Method
hf	-887.50	kJ/mol	Joback Method
hfus	56.10	kJ/mol	Joback Method
hvap	84.05	kJ/mol	Joback Method
log10ws	-7.14		Crippen Method
logp	6.519		Crippen Method
mvol	345.510	ml/mol	McGowan Method
pc	927.24	kPa	Joback Method
rinpol	2545.00		NIST Webbook
rinpol	2545.00		NIST Webbook
tb	874.46	K	Joback Method
tc	1070.59	K	Joback Method
tf	476.53	K	Joback Method
vc	1.347	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1126.53	J/molxK	874.46	Joback Method
cpg	1145.57	J/molxK	907.15	Joback Method
cpg	1163.37	J/molxK	939.84	Joback Method
cpg	1179.96	J/molxK	972.52	Joback Method
cpg	1195.37	J/molxK	1005.21	Joback Method
cpg	1209.63	J/molxK	1037.90	Joback Method
cpg	1222.78	J/molxK	1070.59	Joback Method
dvisc	0.0007041	Paxs	476.53	Joback Method

dvisc	0.0003111	Paxs	542.85	Joback Method
dvisc	0.0001643	Paxs	609.17	Joback Method
dvisc	0.0000983	Paxs	675.50	Joback Method
dvisc	0.0000645	Paxs	741.82	Joback Method
dvisc	0.0000453	Paxs	808.14	Joback Method
dvisc	0.0000336	Paxs	874.46	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U354128&Units=SI

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/51-322-7/Adipic-acid-dodecyl-pent-4-en-2-yl-ester.pdf>

Generated by Cheméo on 2024-04-26 06:54:40.199661764 +0000 UTC m=+16403729.120239079.

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