

Cadina-1(10),4-dien-8«alpha»-ol

Inchi:	InChI=1S/C15H24O/c1-9(2)15-13-7-10(3)5-6-12(13)11(4)8-14(15)16/h7,9,13-16H,5-6,8H
InchiKey:	VAVHKURKIQXFJX-SLTAFYQDSA-N
Formula:	C15H24O
SMILES:	CC1=CC2C(=C(C)CC(O)C2C(C)C)CC1
Mol. weight [g/mol]:	220.35

Physical Properties

Property code	Value	Unit	Source
gf	32.58	kJ/mol	Joback Method
hf	-328.67	kJ/mol	Joback Method
hfus	25.39	kJ/mol	Joback Method
hvap	68.05	kJ/mol	Joback Method
log10ws	-4.25		Crippen Method
logp	3.696		Crippen Method
mcvol	197.760	ml/mol	McGowan Method
pc	2058.62	kPa	Joback Method
rinpol	1678.00		NIST Webbook
rinpol	1680.00		NIST Webbook
ripol	2302.00		NIST Webbook
tb	673.49	K	Joback Method
tc	875.43	K	Joback Method
tf	361.27	K	Joback Method
vc	0.742	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	576.36	J/molxK	673.49	Joback Method
cpg	594.63	J/molxK	707.15	Joback Method
cpg	611.87	J/molxK	740.80	Joback Method
cpg	628.11	J/molxK	774.46	Joback Method
cpg	643.39	J/molxK	808.12	Joback Method
cpg	657.74	J/molxK	841.77	Joback Method
cpg	671.20	J/molxK	875.43	Joback Method

dvisc	0.0034789	Paxs	361.27	Joback Method
dvisc	0.0013534	Paxs	413.31	Joback Method
dvisc	0.0006503	Paxs	465.34	Joback Method
dvisc	0.0003621	Paxs	517.38	Joback Method
dvisc	0.0002244	Paxs	569.42	Joback Method
dvisc	0.0001507	Paxs	621.45	Joback Method
dvisc	0.0001076	Paxs	673.49	Joback Method

Sources

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=R226194&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci9903071
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

Legend

cp_g:	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:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
ripol:	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/56-417-7/Cadina-1-10-4-dien-8-alpha-ol.pdf>

Generated by Cheméo on 2024-04-17 01:28:18.046755444 +0000 UTC m=+15606546.967332759.

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