

(Z)-2-((8R,8aS)-8,8a-Dimethyl-3,4,6,7,8,8a-hexahydro-8,8a-dimethyl-2(1H)-naphthalenyldene)-propanal

Other names:

Propanal,
2-[(8R,8aS)-3,4,6,7,8,8a-hexahydro-8,8a-dimethyl-2(1H)-naphthalenyldene]-propanal, 2-(3,4,6,7,8,8a-hexahydro-8,8a-dimethyl-2(1H)-naphthalenyldene)-propanal, 2-(2Z,8«alpha»,8a«alpha»)-(Z)-Isovalencenal

Inchi: InChI=1S/C15H22O/c1-11(10-16)13-7-8-14-6-4-5-12(2)15(14,3)9-13/h6,10,12H,4-5,7-9H**InchiKey:** SVFQWLHYXVFRHQ-QBFSEMIESA-N**Formula:** C15H22O**SMILES:** CC(C=O)=C1CCC2=CCCC(C)C2(C)C1**Mol. weight [g/mol]:** 218.33**CAS:** 137695-20-6

Physical Properties

Property code	Value	Unit	Source
gf	100.75	kJ/mol	Joback Method
hf	-189.76	kJ/mol	Joback Method
hfus	18.31	kJ/mol	Joback Method
hvap	56.89	kJ/mol	Joback Method
log10ws	-4.39		Crippen Method
logp	4.048		Crippen Method
mcvol	193.460	ml/mol	McGowan Method
pc	2216.62	kPa	Joback Method
rinpol	1812.00		NIST Webbook
rinpol	1812.00		NIST Webbook
tb	632.72	K	Joback Method
tc	861.77	K	Joback Method
tf	356.19	K	Joback Method
vc	0.743	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	527.12	J/molxK	632.72	Joback Method
cpg	547.19	J/molxK	670.90	Joback Method
cpg	566.08	J/molxK	709.07	Joback Method
cpg	583.95	J/molxK	747.25	Joback Method

cpg	600.97	J/mol×K	785.42	Joback Method
cpg	617.33	J/mol×K	823.60	Joback Method
cpg	633.20	J/mol×K	861.77	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C137695206&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

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
mvol:	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/79-050-9/Z-2-8R-8aS-8-8a-Dimethyl-3-4-6-7-8-8a-hexahydronaphthalen-2-1H-ylidene-p>

Generated by Cheméo on 2024-06-15 05:38:16.67315406 +0000 UTC m=+20719145.593731371.

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