

1H-3a,7-Methanoazulene, octahydro-3,6,8,8-tetramethyl-, [3R-(3«alpha»,3a«beta»,6«alpha»,7«beta»,8a«alpha»)-Cedrane]

Inchi: InChI=1S/C15H26/c1-10-7-8-15-9-12(10)14(3,4)13(15)6-5-11(15)2/h10-13H,5-9H2,1-4H3
InchiKey: JJTQQGNEXQKQRF-UHFFFAOYSA-N
Formula: C15H26
SMILES: CC1CCC23CC1C(C)(C)C2CCC3C
Mol. weight [g/mol]: 206.37
CAS: 13567-54-9

Physical Properties

Property code	Value	Unit	Source
gf	199.36	kJ/mol	Joback Method
hf	-177.39	kJ/mol	Joback Method
hfus	15.43	kJ/mol	Joback Method
hvap	45.84	kJ/mol	Joback Method
log10ws	-4.34		Crippen Method
logp	4.495		Crippen Method
mcvol	189.630	ml/mol	McGowan Method
pc	2016.32	kPa	Joback Method
tb	557.83	K	Joback Method
tc	780.06	K	Joback Method
tf	340.67	K	Joback Method
vc	0.724	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	522.52	J/molxK	557.83	Joback Method
cpg	547.59	J/molxK	594.87	Joback Method
cpg	570.94	J/molxK	631.91	Joback Method
cpg	592.86	J/molxK	668.95	Joback Method
cpg	613.64	J/molxK	705.99	Joback Method
cpg	633.54	J/molxK	743.02	Joback Method
cpg	652.86	J/molxK	780.06	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C13567549&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
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvp:	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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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