

(1aR,4S,4aR,7R,7aS,7bS)-1,1,4,7-Tetramethyldeca

Other names:	Epiglobulol
Inchi:	InChI=1S/C15H26O/c1-9-5-6-10-12(9)13-11(14(13,2)3)7-8-15(10,4)16/h9-13,16H,5-8H2,
InchiKey:	AYXPYQRXGNDJFU-IMNVLQEYSA-N
Formula:	C15H26O
SMILES:	CC1CCC2C1C1C(CCC2(C)O)C1(C)C
Mol. weight [g/mol]:	222.37
CAS:	88728-58-9

Physical Properties

Property code	Value	Unit	Source
gf	54.83	kJ/mol	Joback Method
hf	-349.96	kJ/mol	Joback Method
hfus	20.59	kJ/mol	Joback Method
hvap	62.21	kJ/mol	Joback Method
log10ws	-3.71		Crippen Method
logp	3.466		Crippen Method
mcvol	195.500	ml/mol	McGowan Method
pc	2127.56	kPa	Joback Method
rinpol	1587.00		NIST Webbook
rinpol	1584.00		NIST Webbook
rinpol	1581.00		NIST Webbook
rinpol	1608.00		NIST Webbook
rinpol	1597.00		NIST Webbook
rinpol	1584.00		NIST Webbook
rinpol	1597.00		NIST Webbook
rinpol	1564.00		NIST Webbook
rinpol	1564.00		NIST Webbook
rinpol	1543.00		NIST Webbook
rinpol	1540.00		NIST Webbook
rinpol	1608.00		NIST Webbook
rinpol	1590.00		NIST Webbook
rinpol	1585.00		NIST Webbook
rinpol	1580.00		NIST Webbook
rinpol	1580.00		NIST Webbook
rinpol	1585.00		NIST Webbook
rinpol	1564.00		NIST Webbook
rinpol	1564.00		NIST Webbook

rinpol	1585.00		NIST Webbook
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rinpol	1588.00		NIST Webbook
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rinpol	1587.00		NIST Webbook
rinpol	1580.00		NIST Webbook
rinpol	1548.00		NIST Webbook
rinpol	1557.00		NIST Webbook
rinpol	1596.00		NIST Webbook
rinpol	1588.00		NIST Webbook
rinpol	1629.00		NIST Webbook
rinpol	1576.00		NIST Webbook
rinpol	1548.00		NIST Webbook
rinpol	1585.00		NIST Webbook
rinpol	1554.00		NIST Webbook
ripol	2049.00		NIST Webbook
ripol	2049.00		NIST Webbook
ripol	2054.00		NIST Webbook
ripol	2011.00		NIST Webbook
ripol	2025.00		NIST Webbook
ripol	2046.00		NIST Webbook
ripol	2023.00		NIST Webbook
ripol	2025.00		NIST Webbook
ripol	2033.00		NIST Webbook
ripol	2054.00		NIST Webbook
ripol	2016.00		NIST Webbook
tb	645.34	K	Joback Method
tc	851.43	K	Joback Method
tf	397.25	K	Joback Method
vc	0.742	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	600.58	J/mol×K	645.34	Joback Method
cpg	621.09	J/mol×K	679.69	Joback Method
cpg	640.66	J/mol×K	714.04	Joback Method
cpg	659.50	J/mol×K	748.38	Joback Method
cpg	677.83	J/mol×K	782.73	Joback Method
cpg	695.86	J/mol×K	817.08	Joback Method
cpg	713.82	J/mol×K	851.43	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=C88728589&Units=SI

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
ripol:	Non-polar retention indices
ripol:	Polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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