

Amorpha-4,9-dien-2«alpha»-ol

Inchi:

InchiKey:

Formula:

SMILES:

Mol. weight [g/mol]:

InChI=1S/C15H24O/c1-9(2)12-6-5-11(4)15-13(12)7-10(3)8-14(15)16/h5,7,9,12-16H,6,8H

STLBTFMCSEXAQ-XGUBFFRZSA-N

C15H24O

CC1=CC2C(C(C)C)CC=C(C)C2C(O)C1

220.35

Physical Properties

Property code	Value	Unit	Source
gf	34.50	kJ/mol	Joback Method
hf	-337.54	kJ/mol	Joback Method
hfus	26.85	kJ/mol	Joback Method
hvap	67.08	kJ/mol	Joback Method
log10ws	-4.01		Crippen Method
logp	3.552		Crippen Method
mcvol	197.760	ml/mol	McGowan Method
pc	2025.41	kPa	Joback Method
rinpol	1713.00		NIST Webbook
rinpol	1713.00		NIST Webbook
tb	663.84	K	Joback Method
tc	864.81	K	Joback Method
tf	344.51	K	Joback Method
vc	0.741	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	578.38	J/mol×K	663.84	Joback Method
cpg	597.29	J/mol×K	697.34	Joback Method
cpg	615.12	J/mol×K	730.83	Joback Method
cpg	631.91	J/mol×K	764.33	Joback Method
cpg	647.70	J/mol×K	797.82	Joback Method
cpg	662.51	J/mol×K	831.32	Joback Method
cpg	676.39	J/mol×K	864.81	Joback Method
dvisc	0.0043769	Pa×s	344.51	Joback Method

dvisc	0.0016494	Pa×s	397.73	Joback Method
dvisc	0.0007826	Pa×s	450.95	Joback Method
dvisc	0.0004346	Pa×s	504.17	Joback Method
dvisc	0.0002700	Pa×s	557.40	Joback Method
dvisc	0.0001823	Pa×s	610.62	Joback Method
dvisc	0.0001311	Pa×s	663.84	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=R639926&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cp_g:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
g_f:	Standard Gibbs free energy of formation
h_f:	Enthalpy of formation at standard conditions
h_{fus}:	Enthalpy of fusion at standard conditions
h_{vap}:	Enthalpy of vaporization at standard conditions
log_{10ws}:	Log10 of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
p_c:	Critical Pressure
rin_{pol}:	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/71-903-0/Amorpha-4-9-dien-2-alpha-ol.pdf>

Generated by Cheméo on 2025-12-24 12:06:01.518918005 +0000 UTC m=+6326159.048958663.

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