

Tris(2-methylallyl)amine

Other names:	(CH ₂ =C(CH ₃)CH ₂) ₃ N 2-Propen-1-amine, 2-methyl-N,N-bis(2-methyl-2-propenyl)-
Inchi:	InChI=1S/C12H21N/c1-10(2)7-13(8-11(3)4)9-12(5)6/h1,3,5,7-9H2,2,4,6H3
InchiKey:	FXBJYRVIFGLPBC-UHFFFAOYSA-N
Formula:	C ₁₂ H ₂₁ N
SMILES:	C=C(C)CN(CC(=C)C)CC(=C)C
Mol. weight [g/mol]:	179.30
CAS:	6321-40-0

Physical Properties

Property code	Value	Unit	Source
affp	980.20	kJ/mol	NIST Webbook
basg	949.40	kJ/mol	NIST Webbook
gf	398.81	kJ/mol	Joback Method
hf	123.44	kJ/mol	Joback Method
hfus	22.09	kJ/mol	Joback Method
hvap	42.58	kJ/mol	Joback Method
ie	7.80	eV	NIST Webbook
log10ws	-2.98		Crippen Method
logp	3.017		Crippen Method
mcvol	177.020	ml/mol	McGowan Method
pc	2009.10	kPa	Joback Method
tb	476.08	K	Joback Method
tc	653.74	K	Joback Method
tf	210.31	K	Joback Method
vc	0.671	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	387.32	J/mol×K	476.08	Joback Method
cpg	404.17	J/mol×K	505.69	Joback Method
cpg	420.19	J/mol×K	535.30	Joback Method
cpg	435.41	J/mol×K	564.91	Joback Method

cpg	449.88	J/mol×K	594.52	Joback Method
cpg	463.62	J/mol×K	624.13	Joback Method
cpg	476.67	J/mol×K	653.74	Joback Method

Pressure Dependent Properties

Property code	Value	Unit	Pressure [kPa]	Source
tbrp	357.20	K	2.00	NIST Webbook

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C6321400&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
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772

Legend

affp:	Proton affinity
basg:	Gas basicity
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
ie:	Ionization energy
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
tbrp:	Boiling point at reduced pressure
tc:	Critical Temperature
tf:	Normal melting (fusion) point

vc: Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/79-398-4/Tris-2-methylallyl-amine.pdf>

Generated by Cheméo on 2024-04-30 18:40:37.809763633 +0000 UTC m=+16791686.730340944.

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