

(2E,4E,14E)-N-Isobutylicos-2,4,14-trienamide

Inchi: InChI=1S/C24H43NO/c1-4-5-6-7-8-9-10-11-12-13-14-15-16-17-18-19-20-21-24(26)25-22
InchiKey: GQCWFFNZERNJJC-DIIODVDPSA-N
Formula: C24H43NO
SMILES: CCCCCC=CCCCCCCCC=CC=CC(O)=NCC(C)C
Mol. weight [g/mol]: 361.60
CAS: 943546-21-2

Physical Properties

Property code	Value	Unit	Source
hf	-272.11	kJ/mol	Joback Method
hvap	88.58	kJ/mol	Joback Method
log10ws	-8.17		Crippen Method
logp	7.969		Crippen Method
mcvol	347.670	ml/mol	McGowan Method
pc	882.62	kPa	Joback Method
rinpol	2985.00		NIST Webbook
rinpol	2985.00		NIST Webbook
tb	929.30	K	Joback Method
tc	1137.88	K	Joback Method

Sources

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>
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C943546212&Units=SI>

Legend

hf: Enthalpy of formation at standard conditions

h_{vap}:	Enthalpy of vaporization at standard conditions
log₁₀w_s:	Log ₁₀ of Water solubility in mol/l
log_p:	Octanol/Water partition coefficient
mc_{vol}:	McGowan's characteristic volume
p_c:	Critical Pressure
r_{inpol}:	Non-polar retention indices
t_b:	Normal Boiling Point Temperature
t_c:	Critical Temperature

Latest version available from:

<https://www.cheméo.com/cid/95-645-1/2E-4E-14E-N-Isobutylicos-2-4-14-trienamide.pdf>

Generated by Cheméo on 2024-05-12 18:44:02.272453254 +0000 UTC m=+17828691.193030567.

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