

progesterone, MO-TMS, syn

Inchi: InChI=1S/C23H36N2O2/c1-15(24-26-4)19-8-9-20-18-7-6-16-14-17(25-27-5)10-12-22(16,
InchiKey: WCBFSBXYZBTGBQ-JYVPLEQZSA-N
Formula: C23H36N2O2
SMILES: CON=C1C=C2CCC3C(CCC4(C)C(C(C)=NOC)CCC34)C2(C)CC1
Mol. weight [g/mol]: 372.54

Physical Properties

Property code	Value	Unit	Source
hf	-372.52	kJ/mol	Joback Method
hvap	77.70	kJ/mol	Joback Method
log10ws	-5.92		Crippen Method
logp	5.590		Crippen Method
mcvol	310.290	ml/mol	McGowan Method
pc	1112.59	kPa	Joback Method
rinpol	2907.00		NIST Webbook
tb	969.79	K	Joback Method
tc	1217.22	K	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=R395869&Units=SI>

Legend

hf: Enthalpy of formation 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
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature

Latest version available from:

<https://www.cheméo.com/cid/70-499-1/progesterone-MO-TMS-syn.pdf>

Generated by Cheméo on 2024-04-27 15:46:07.164241038 +0000 UTC m=+16522016.084818353.

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