p-methoxybenzylidene-nonyl-amine

Inchi: InChI=1S/C17H27NO/c1-3-4-5-6-7-8-9-14-18-15-16-10-12-17(19-2)13-11-16/h10-13,15H
InchiKey: HUFQEOSXDHSMIW-UHFFFAOYSA-N
Formula: C17H27NO
SMILES: CCCCCCCCCN=Cc1ccc(OC)cc1
Mol. weight [g/mol]: 261.40

Physical Properties

Property code	Value	Unit	Source
hf	-219.15	kJ/mol	Joback Method
hvap	62.10	kJ/mol	Joback Method
log10ws	-5.01		Crippen Method
logp	4.865		Crippen Method
mcvol	238.180	ml/mol	McGowan Method
pc	1417.57	kPa	Joback Method
rinpol	2130.00		NIST Webbook
tb	719.12	K	Joback Method
tc	919.56	K	Joback Method

Sources

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=R160148&Units=SI
Crippen Method: http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws

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