

# p-Butoxybenzylidene p-hexylaniline

**Inchi:** InChI=1S/C23H31NO/c1-3-5-7-8-9-20-10-14-22(15-11-20)24-19-21-12-16-23(17-13-21)2  
**InchiKey:** OFIBDXCRLAWBQU-UHFFFAOYSA-N  
**Formula:** C23H31NO  
**SMILES:** CCCCCc1ccc(N=Cc2ccc(OCCCC)cc2)cc1  
**Mol. weight [g/mol]:** 337.50  
**CAS:** 39777-12-3

## Physical Properties

Property code	Value	Unit	Source
hf	-117.93	kJ/mol	Joback Method
hvap	78.39	kJ/mol	Joback Method
log10ws	-7.19		Crippen Method
logp	6.739		Crippen Method
mvol	298.960	ml/mol	McGowan Method
pc	1176.05	kPa	Joback Method
tb	888.06	K	Joback Method
tc	1108.47	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](https://www.chemeo.com/doc/models/crippen_log10ws)  
**Joback Method:** [https://en.wikipedia.org/wiki/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=C39777123&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

<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

Latest version available from:

<https://www.chemeo.com/cid/51-560-3/p-Butoxybenzylidene-p-hexylaniline.pdf>

Generated by Cheméo on 2024-02-24 11:47:04.717511836 +0000 UTC m=+11064473.638089151.

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