

# p-methylbenzylidene-(4-ethoxyphenyl)-amine

**Inchi:** InChI=1S/C16H17NO/c1-3-18-16-10-8-15(9-11-16)17-12-14-6-4-13(2)5-7-14/h4-12H,3H2  
**InchiKey:** YLVDCPUKGYDWBK-SFQUDFHCSA-N  
**Formula:** C16H17NO  
**SMILES:** CCOc1ccc(N=Cc2ccc(C)cc2)cc1  
**Mol. weight [g/mol]:** 239.31

## Physical Properties

Property code	Value	Unit	Source
hf	26.55	kJ/mol	Joback Method
hvap	62.81	kJ/mol	Joback Method
log10ws	-4.35		Crippen Method
logp	4.144		Crippen Method
mcvol	200.330	ml/mol	McGowan Method
pc	1985.89	kPa	Joback Method
rinpol	2212.00		NIST Webbook
rinpol	2212.00		NIST Webbook
tb	727.90	K	Joback Method
tc	969.82	K	Joback Method

## Sources

**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=R160282&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.chemeo.com/doc/models/crippen\\_log10ws](https://www.chemeo.com/doc/models/crippen_log10ws)

## Legend

**hf:** Enthalpy of formation at standard conditions  
**hvap:** Enthalpy of vaporization at standard conditions

<b>log10ws:</b>	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>rinpol:</b>	Non-polar retention indices
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature

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