

Methacrylic acid, 4-biphenyl ester

Inchi:	InChI=1S/C16H14O2/c1-12(2)16(17)18-15-10-8-14(9-11-15)13-6-4-3-5-7-13/h3-11H,1H2
InchiKey:	HEJFLAMVALNBRR-UHFFFAOYSA-N
Formula:	C16H14O2
SMILES:	<chem>C=C(C)C(=O)Oc1ccc(-c2ccccc2)cc1</chem>
Mol. weight [g/mol]:	238.28

Physical Properties

Property code	Value	Unit	Source
gf	144.40	kJ/mol	Joback Method
hf	-41.14	kJ/mol	Joback Method
hfus	25.09	kJ/mol	Joback Method
hvap	64.99	kJ/mol	Joback Method
log10ws	-5.08		Crippen Method
logp	3.835		Crippen Method
mvol	191.920	ml/mol	McGowan Method
pc	2465.36	kPa	Joback Method
rinpol	2025.00		NIST Webbook
rinpol	2025.00		NIST Webbook
tb	696.67	K	Joback Method
tc	939.18	K	Joback Method
tf	391.88	K	Joback Method
vc	0.722	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	494.08	J/molxK	696.67	Joback Method
cpg	509.67	J/molxK	737.09	Joback Method
cpg	524.01	J/molxK	777.51	Joback Method
cpg	537.16	J/molxK	817.93	Joback Method
cpg	549.19	J/molxK	858.34	Joback Method
cpg	560.17	J/molxK	898.76	Joback Method
cpg	570.15	J/molxK	939.18	Joback Method

Sources

NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U360695&Units=SI
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

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvp:	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
rinp:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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