

# Adipic acid, 4-biphenyl ethyl ester

<b>Inchi:</b>	InChI=1S/C20H22O4/c1-2-23-19(21)10-6-7-11-20(22)24-18-14-12-17(13-15-18)16-8-4-3
<b>InchiKey:</b>	KQBDEIRQOFZBMJ-UHFFFAOYSA-N
<b>Formula:</b>	C20H22O4
<b>SMILES:</b>	CCOC(=O)CCCCC(=O)Oc1ccc(-c2ccccc2)cc1
<b>Mol. weight [g/mol]:</b>	326.39

## Physical Properties

Property code	Value	Unit	Source
gf	-135.13	kJ/mol	Joback Method
hf	-484.14	kJ/mol	Joback Method
hfus	40.82	kJ/mol	Joback Method
hvap	83.64	kJ/mol	Joback Method
log10ws	-5.76		Crippen Method
logp	4.382		Crippen Method
mcvol	260.020	ml/mol	McGowan Method
pc	1737.56	kPa	Joback Method
rinpol	2688.00		NIST Webbook
rinpol	2688.00		NIST Webbook
tb	867.92	K	Joback Method
tc	1091.70	K	Joback Method
tf	524.84	K	Joback Method
vc	0.988	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	781.49	J/molxK	867.92	Joback Method
cpg	795.69	J/molxK	905.22	Joback Method
cpg	808.60	J/molxK	942.51	Joback Method
cpg	820.28	J/molxK	979.81	Joback Method
cpg	830.75	J/molxK	1017.10	Joback Method
cpg	840.06	J/molxK	1054.40	Joback Method
cpg	848.25	J/molxK	1091.70	Joback Method
dvisc	0.0005161	Paxs	524.84	Joback Method

dvisc	0.0002974	Paxs	582.02	Joback Method
dvisc	0.0001891	Paxs	639.20	Joback Method
dvisc	0.0001295	Paxs	696.38	Joback Method
dvisc	0.0000939	Paxs	753.56	Joback Method
dvisc	0.0000713	Paxs	810.74	Joback Method
dvisc	0.0000561	Paxs	867.92	Joback Method

## Sources

<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=U353920&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=U353920&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	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
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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