

4-Ethylbenzoic acid, 2-adamantyl ester

Inchi:	InChI=1S/C19H24O2/c1-2-12-3-5-15(6-4-12)19(20)21-18-16-8-13-7-14(10-16)11-17(18)9
InchiKey:	ZIIYEEGFRAAVLV-UHFFFAOYSA-N
Formula:	C19H24O2
SMILES:	CCc1ccc(C(=O)OC2C3CC4CC(C3)CC2C4)cc1
Mol. weight [g/mol]:	284.39

Physical Properties

Property code	Value	Unit	Source
gf	132.69	kJ/mol	Joback Method
hf	-283.67	kJ/mol	Joback Method
hfus	35.85	kJ/mol	Joback Method
hvap	69.28	kJ/mol	Joback Method
log10ws	-5.11		Crippen Method
logp	4.231		Crippen Method
mcvol	229.670	ml/mol	McGowan Method
pc	1840.41	kPa	Joback Method
tb	757.22	K	Joback Method
tc	986.75	K	Joback Method
tf	456.81	K	Joback Method
vc	0.876	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	731.10	J/molxK	757.22	Joback Method
cpg	820.35	J/molxK	948.49	Joback Method
cpg	804.83	J/molxK	910.24	Joback Method
cpg	788.26	J/molxK	871.98	Joback Method
cpg	770.54	J/molxK	833.73	Joback Method
cpg	751.52	J/molxK	795.47	Joback Method
cpg	834.95	J/molxK	986.75	Joback Method
dvisc	0.0022310	Paxs	757.22	Joback Method
dvisc	0.0023642	Paxs	707.15	Joback Method
dvisc	0.0025277	Paxs	657.08	Joback Method

dvisc	0.0027324	Paxs	607.01	Joback Method
dvisc	0.0029953	Paxs	556.95	Joback Method
dvisc	0.0033437	Paxs	506.88	Joback Method
dvisc	0.0038237	Paxs	456.81	Joback Method

Sources

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
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U292198&Units=SI

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvac:	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
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

Latest version available from:

<https://www.chemeo.com/cid/89-502-6/4-Ethylbenzoic-acid-2-adamantyl-ester.pdf>

Generated by Cheméo on 2024-04-20 12:09:46.984228799 +0000 UTC m=+15904235.904806111.

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