Pentacyclo[13.3.2.26,10.13,18.19,12]tetracosa-1,3(

Other names: Pentacyclo[13.3.2.2
Inchi: InChI=1S/C24H20/c1-2-18-7-11-23-15-20(16-24(23)12-8-18)4-3-19-13-21-9-5-17(1)6-10-
InchiKey: SOJCCVPXRNRYQH-UHFFFAOYSA-N
Formula: C24H20
SMILES: c1cc2cc3cc-2ccc1CCc1ccc2cc(cc-2cc1)CC3
Mol. weight [g/mol]: 308.42

Physical Properties

Property code	Value	Unit	Source
hf	373.52	kJ/mol	Cheméo Relay (1.0)
hfus	33.44	kJ/mol	Joback Method
ie	6.60	eV	NIST Webbook
logp	5.780		Crippen Method
mcvol	251.720	ml/mol	McGowan Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	755.43	J/mol×K	875.44	Joback Method
cpg	772.92	J/mol×K	920.50	Joback Method
cpg	789.45	J/mol×K	965.56	Joback Method
cpg	805.27	J/mol×K	1010.62	Joback Method
cpg	820.66	J/mol×K	1055.68	Joback Method
cpg	835.89	J/mol×K	1100.74	Joback Method
cpg	851.22	J/mol×K	1145.80	Joback Method
dvisc	0.0014536	Pa×s	547.22	Joback Method
dvisc	0.0010812	Pa×s	601.92	Joback Method
dvisc	0.0008448	Pa×s	656.63	Joback Method
dvisc	0.0006857	Pa×s	711.33	Joback Method
dvisc	0.0005733	Pa×s	766.03	Joback Method
dvisc	0.0004910	Pa×s	820.74	Joback Method
dvisc	0.0004287	Pa×s	875.44	Joback Method

Sources

McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Cheméo Relay (1.0):	
Cheméo Relay (1.0, beta):	
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=C73608523&Units=SI
Joback Method:	https://en.wikipedia.org/wiki/Joback_method

Legend

cpg:	Ideal gas heat capacity
dvisc:	Dynamic viscosity
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
ie:	Ionization energy
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume