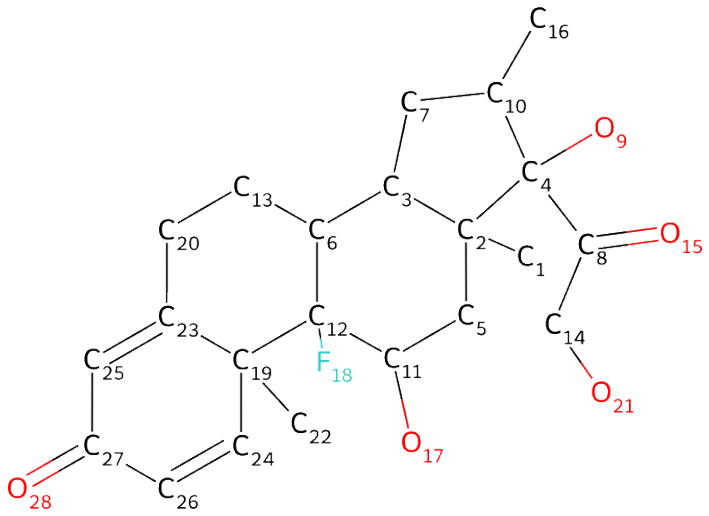
C-oxidation BDE Energy Report for: 107_dexamethasone-out

This report covers the results for bond dissociation enthalpies (BDE) and solvent accessible surface area (SASA) calculations performed for 107_dexamethasone-out. Oxidation propensity is established using C-H BDE. The lower the C-H BDE values the higher the propensity for C-oxidation. Details for the density functional theory (DFT) calculations and overall workflow are explained at the end of this document.

BDE and SASA



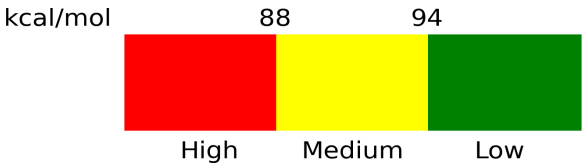
Atom	BDE (kcal/mol)	Propensity	SASA (Å²)
C1	100.05	Low	3.86
C3	87.23	High	0.00
C5	97.89	Low	0.26
C6	96.59	Low	0.00
C7	96.69	Low	0.25
C10	91.36	Moderate	0.00
C11	90.23	Moderate	0.00
C13	97.8	Low	0.41
C14	75.33	High	1.94
C16	99.12	Low	5.33
C20	82.62	High	1.30
C22	99.77	Low	4.33
C24	106.7	Low	1.74

C25	112.36	Low	9.06
C26	112.38	Low	12.45

Missing Sites:

None

Risk Scale:



Calculation Details

Conformational search calculations were performed only for the base ground state molecule. The lowest energy conformer was selected to generate radicals and run optimization DFT calculations. DFT calculations were performed using Gaussian with B3LYP level of theory and 6-31G(d,p) basis set. The BDE protocol was adapted from: *Lienard, P., Gavartin, J., Boccardi, G., & Meunier, M. (2015). Predicting drug substances autoxidation. Pharmaceutical research, 32, 300-310.*