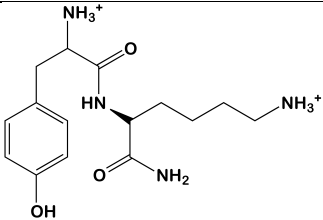
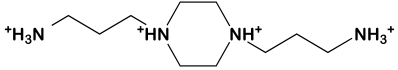
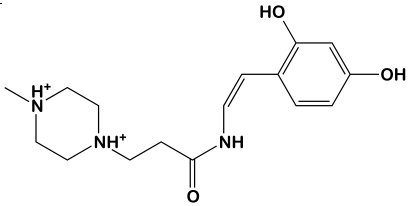
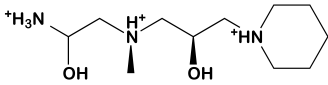
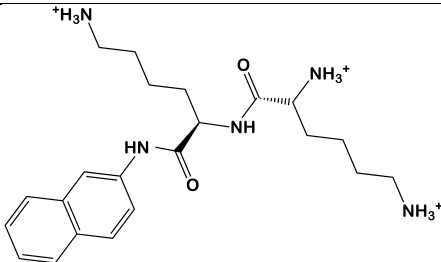


**Additional file 1** Additional compounds identified from virtual screening that were docked and tested *in vitro* against PfSpdS.

Compound	Structure	Best Fit Value/number of PhFs	Docking scores (kcal/mol)	% Inhibition (at 100 $\mu$ M) <sup>a</sup>
<b>DPM2</b>				
3		0.91/4	-12.5	No Inhibition
4		0.15/4	-9.8	No Inhibition
<b>DPM3</b>				
5		0.10/5	-9.0	No Inhibition
6		0.18/4	-9.7	No Inhibition
<b>DPM4</b>				
7		0.14/5	-10.5	No Inhibition

<sup>a</sup> Results represent inhibition of PfSpdS activity compared to untreated enzyme of three independent experiments performed in duplicate.