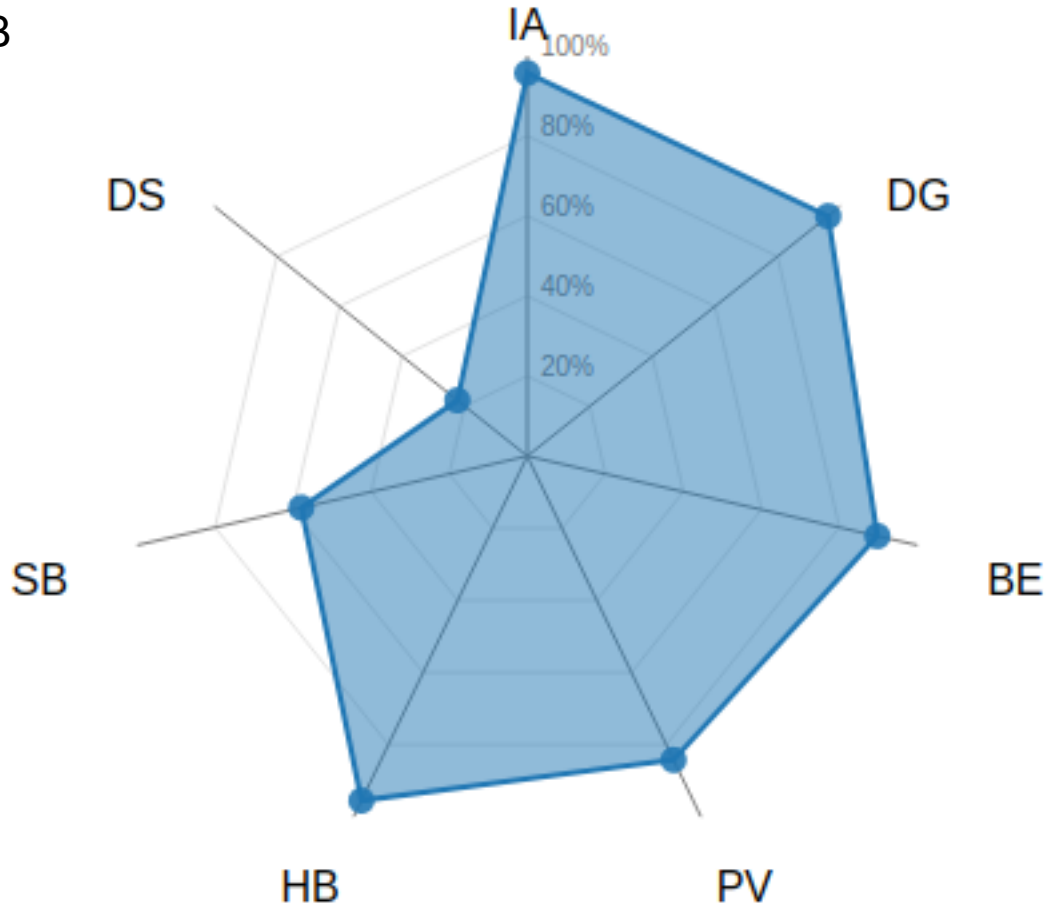
Interaction radar

chains A-B

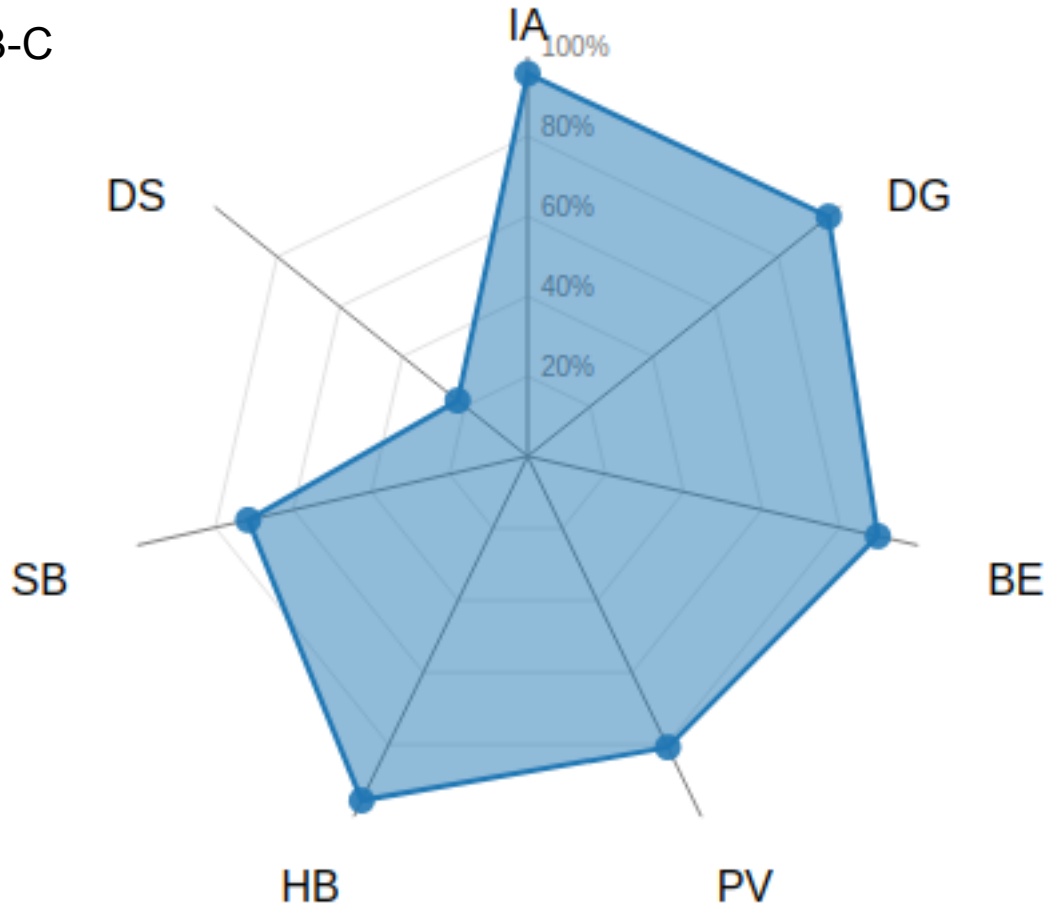


Interface parameters

IA	:	Interface area, Å^2	8443
DG	:	Solvation Energy, kcal/mol	-103.8
BE	:	Total Binding Energy, kcal/mol	-124.5
PV	:	Hydrophobic P-value	0.03078
HB	:	Number of Hydrogen Bonds	40
SB	:	Number of Salt Bridges	8
DS	:	Number of Disulphide Bonds	0

Interaction radar

chains B-C

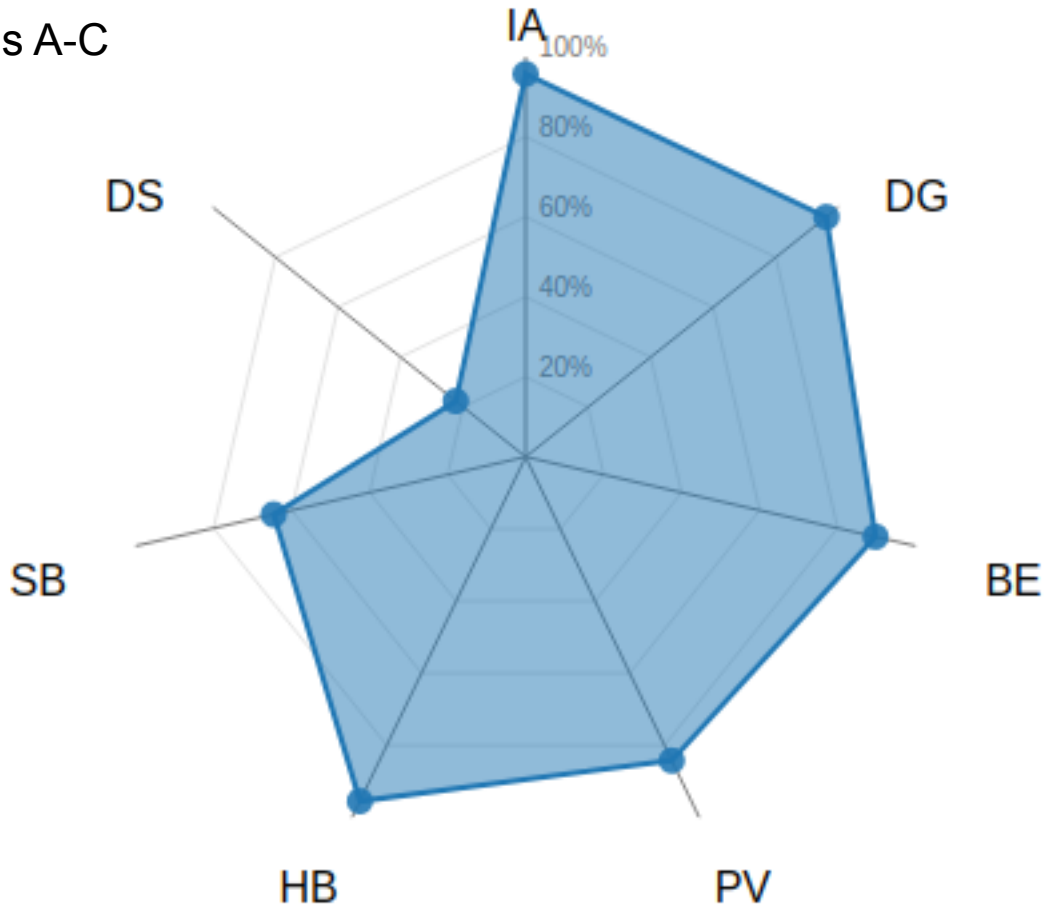


Interface parameters

IA	:	Interface area, Å^2	8415
DG	:	Solvation Energy, kcal/mol	-103.4
BE	:	Total Binding Energy, kcal/mol	-124.4
PV	:	Hydrophobic P-value	0.04092
HB	:	Number of Hydrogen Bonds	38
SB	:	Number of Salt Bridges	11
DS	:	Number of Disulphide Bonds	0

Interaction radar

chains A-C



Interface parameters

IA	:	Interface area, Å²	8421
DG	:	Solvation Energy, kcal/mol	-103.3
BE	:	Total Binding Energy, kcal/mol	-123.9
PV	:	Hydrophobic P-value	0.03349
HB	:	Number of Hydrogen Bonds	38
SB	:	Number of Salt Bridges	10
DS	:	Number of Disulphide Bonds	0