



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 17, 2024 – 05:41 PM EDT

PDB ID : 3DTD  
Title : Crystal structure of invasion associated protein b from bartonella henselae  
Authors : Patskovsky, Y.; Ozyurt, S.; Freeman, J.; Slocombe, A.; Groshong, C.; Koss, J.; Smith, D.; Wasserman, S.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2008-07-14  
Resolution : 2.35 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

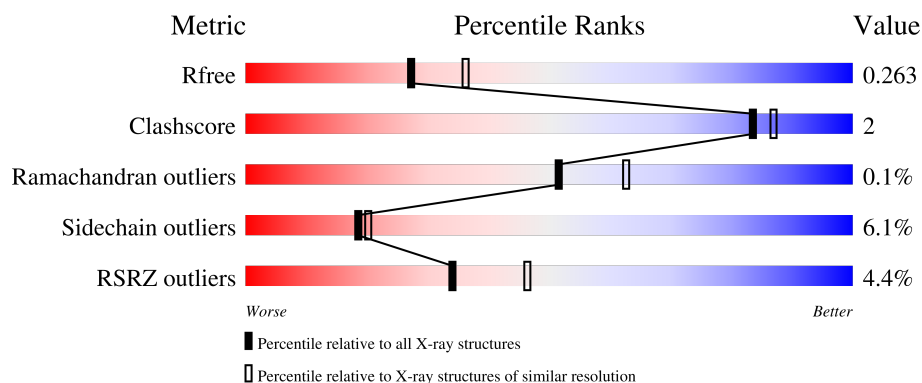
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	175	<div> <div>5%</div> <div> <div></div> <div>71%</div> <div>10%</div> <div>•</div> <div>19%</div> </div> </div>
1	B	175	<div> <div>%</div> <div> <div></div> <div>75%</div> <div>7%</div> <div>•</div> <div>17%</div> </div> </div>
1	C	175	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>9%</div> <div>•</div> <div>16%</div> </div> </div>
1	D	175	<div> <div>4%</div> <div> <div></div> <div>71%</div> <div>11%</div> <div>•</div> <div>17%</div> </div> </div>
1	E	175	<div> <div>7%</div> <div> <div></div> <div>77%</div> <div>5%</div> <div>•</div> <div>18%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	175	
1	G	175	
1	H	175	
1	I	175	
1	J	175	
1	K	175	
1	L	175	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	D	1	-	-	-	X
2	GOL	F	1	-	-	X	-
2	GOL	G	1	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13760 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Invasion-associated protein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	142	Total	C	N	O	S	0	0	0
			1074	685	188	194	7			
1	B	146	Total	C	N	O	S	0	2	0
			1113	708	197	201	7			
1	C	147	Total	C	N	O	S	0	1	0
			1113	708	196	202	7			
1	D	145	Total	C	N	O	S	0	2	0
			1109	706	196	200	7			
1	E	143	Total	C	N	O	S	0	1	0
			1091	696	193	195	7			
1	F	142	Total	C	N	O	S	0	2	0
			1090	695	194	194	7			
1	G	143	Total	C	N	O	S	0	2	0
			1097	699	193	198	7			
1	H	146	Total	C	N	O	S	0	2	0
			1115	709	197	202	7			
1	I	145	Total	C	N	O	S	0	2	0
			1108	707	195	199	7			
1	J	145	Total	C	N	O	S	0	2	0
			1107	705	195	200	7			
1	K	145	Total	C	N	O	S	0	2	0
			1112	708	198	199	7			
1	L	145	Total	C	N	O	S	0	2	0
			1107	705	195	200	7			

There are 132 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	MET	-	expression tag	UNP Q6G4Y3
A	20	SER	-	expression tag	UNP Q6G4Y3
A	21	LEU	-	expression tag	UNP Q6G4Y3
A	186	GLU	-	expression tag	UNP Q6G4Y3
A	187	GLY	-	expression tag	UNP Q6G4Y3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	188	HIS	-	expression tag	UNP Q6G4Y3
A	189	HIS	-	expression tag	UNP Q6G4Y3
A	190	HIS	-	expression tag	UNP Q6G4Y3
A	191	HIS	-	expression tag	UNP Q6G4Y3
A	192	HIS	-	expression tag	UNP Q6G4Y3
A	193	HIS	-	expression tag	UNP Q6G4Y3
B	19	MET	-	expression tag	UNP Q6G4Y3
B	20	SER	-	expression tag	UNP Q6G4Y3
B	21	LEU	-	expression tag	UNP Q6G4Y3
B	186	GLU	-	expression tag	UNP Q6G4Y3
B	187	GLY	-	expression tag	UNP Q6G4Y3
B	188	HIS	-	expression tag	UNP Q6G4Y3
B	189	HIS	-	expression tag	UNP Q6G4Y3
B	190	HIS	-	expression tag	UNP Q6G4Y3
B	191	HIS	-	expression tag	UNP Q6G4Y3
B	192	HIS	-	expression tag	UNP Q6G4Y3
B	193	HIS	-	expression tag	UNP Q6G4Y3
C	19	MET	-	expression tag	UNP Q6G4Y3
C	20	SER	-	expression tag	UNP Q6G4Y3
C	21	LEU	-	expression tag	UNP Q6G4Y3
C	186	GLU	-	expression tag	UNP Q6G4Y3
C	187	GLY	-	expression tag	UNP Q6G4Y3
C	188	HIS	-	expression tag	UNP Q6G4Y3
C	189	HIS	-	expression tag	UNP Q6G4Y3
C	190	HIS	-	expression tag	UNP Q6G4Y3
C	191	HIS	-	expression tag	UNP Q6G4Y3
C	192	HIS	-	expression tag	UNP Q6G4Y3
C	193	HIS	-	expression tag	UNP Q6G4Y3
D	19	MET	-	expression tag	UNP Q6G4Y3
D	20	SER	-	expression tag	UNP Q6G4Y3
D	21	LEU	-	expression tag	UNP Q6G4Y3
D	186	GLU	-	expression tag	UNP Q6G4Y3
D	187	GLY	-	expression tag	UNP Q6G4Y3
D	188	HIS	-	expression tag	UNP Q6G4Y3
D	189	HIS	-	expression tag	UNP Q6G4Y3
D	190	HIS	-	expression tag	UNP Q6G4Y3
D	191	HIS	-	expression tag	UNP Q6G4Y3
D	192	HIS	-	expression tag	UNP Q6G4Y3
D	193	HIS	-	expression tag	UNP Q6G4Y3
E	19	MET	-	expression tag	UNP Q6G4Y3
E	20	SER	-	expression tag	UNP Q6G4Y3
E	21	LEU	-	expression tag	UNP Q6G4Y3

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Chain	Residue	Modelled	Actual	Comment	Reference
E	186	GLU	-	expression tag	UNP Q6G4Y3
E	187	GLY	-	expression tag	UNP Q6G4Y3
E	188	HIS	-	expression tag	UNP Q6G4Y3
E	189	HIS	-	expression tag	UNP Q6G4Y3
E	190	HIS	-	expression tag	UNP Q6G4Y3
E	191	HIS	-	expression tag	UNP Q6G4Y3
E	192	HIS	-	expression tag	UNP Q6G4Y3
E	193	HIS	-	expression tag	UNP Q6G4Y3
F	19	MET	-	expression tag	UNP Q6G4Y3
F	20	SER	-	expression tag	UNP Q6G4Y3
F	21	LEU	-	expression tag	UNP Q6G4Y3
F	186	GLU	-	expression tag	UNP Q6G4Y3
F	187	GLY	-	expression tag	UNP Q6G4Y3
F	188	HIS	-	expression tag	UNP Q6G4Y3
F	189	HIS	-	expression tag	UNP Q6G4Y3
F	190	HIS	-	expression tag	UNP Q6G4Y3
F	191	HIS	-	expression tag	UNP Q6G4Y3
F	192	HIS	-	expression tag	UNP Q6G4Y3
F	193	HIS	-	expression tag	UNP Q6G4Y3
G	19	MET	-	expression tag	UNP Q6G4Y3
G	20	SER	-	expression tag	UNP Q6G4Y3
G	21	LEU	-	expression tag	UNP Q6G4Y3
G	186	GLU	-	expression tag	UNP Q6G4Y3
G	187	GLY	-	expression tag	UNP Q6G4Y3
G	188	HIS	-	expression tag	UNP Q6G4Y3
G	189	HIS	-	expression tag	UNP Q6G4Y3
G	190	HIS	-	expression tag	UNP Q6G4Y3
G	191	HIS	-	expression tag	UNP Q6G4Y3
G	192	HIS	-	expression tag	UNP Q6G4Y3
G	193	HIS	-	expression tag	UNP Q6G4Y3
H	19	MET	-	expression tag	UNP Q6G4Y3
H	20	SER	-	expression tag	UNP Q6G4Y3
H	21	LEU	-	expression tag	UNP Q6G4Y3
H	186	GLU	-	expression tag	UNP Q6G4Y3
H	187	GLY	-	expression tag	UNP Q6G4Y3
H	188	HIS	-	expression tag	UNP Q6G4Y3
H	189	HIS	-	expression tag	UNP Q6G4Y3
H	190	HIS	-	expression tag	UNP Q6G4Y3
H	191	HIS	-	expression tag	UNP Q6G4Y3
H	192	HIS	-	expression tag	UNP Q6G4Y3
H	193	HIS	-	expression tag	UNP Q6G4Y3
I	19	MET	-	expression tag	UNP Q6G4Y3

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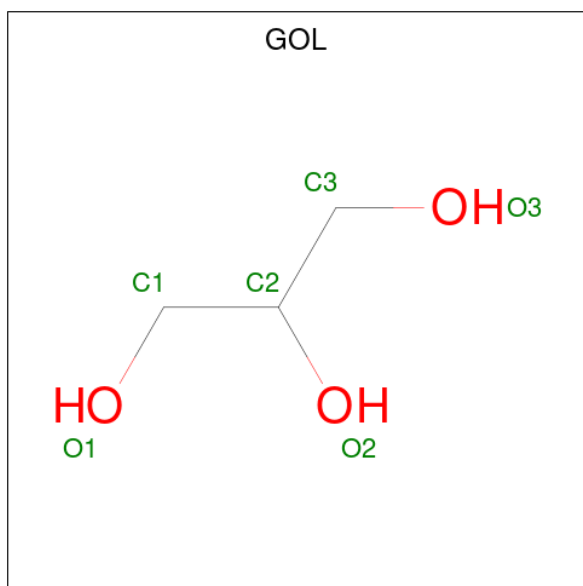
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I	21	LEU	-	expression tag	UNP Q6G4Y3
I	186	GLU	-	expression tag	UNP Q6G4Y3
I	187	GLY	-	expression tag	UNP Q6G4Y3
I	188	HIS	-	expression tag	UNP Q6G4Y3
I	189	HIS	-	expression tag	UNP Q6G4Y3
I	190	HIS	-	expression tag	UNP Q6G4Y3
I	191	HIS	-	expression tag	UNP Q6G4Y3
I	192	HIS	-	expression tag	UNP Q6G4Y3
I	193	HIS	-	expression tag	UNP Q6G4Y3
J	19	MET	-	expression tag	UNP Q6G4Y3
J	20	SER	-	expression tag	UNP Q6G4Y3
J	21	LEU	-	expression tag	UNP Q6G4Y3
J	186	GLU	-	expression tag	UNP Q6G4Y3
J	187	GLY	-	expression tag	UNP Q6G4Y3
J	188	HIS	-	expression tag	UNP Q6G4Y3
J	189	HIS	-	expression tag	UNP Q6G4Y3
J	190	HIS	-	expression tag	UNP Q6G4Y3
J	191	HIS	-	expression tag	UNP Q6G4Y3
J	192	HIS	-	expression tag	UNP Q6G4Y3
J	193	HIS	-	expression tag	UNP Q6G4Y3
K	19	MET	-	expression tag	UNP Q6G4Y3
K	20	SER	-	expression tag	UNP Q6G4Y3
K	21	LEU	-	expression tag	UNP Q6G4Y3
K	186	GLU	-	expression tag	UNP Q6G4Y3
K	187	GLY	-	expression tag	UNP Q6G4Y3
K	188	HIS	-	expression tag	UNP Q6G4Y3
K	189	HIS	-	expression tag	UNP Q6G4Y3
K	190	HIS	-	expression tag	UNP Q6G4Y3
K	191	HIS	-	expression tag	UNP Q6G4Y3
K	192	HIS	-	expression tag	UNP Q6G4Y3
K	193	HIS	-	expression tag	UNP Q6G4Y3
L	19	MET	-	expression tag	UNP Q6G4Y3
L	20	SER	-	expression tag	UNP Q6G4Y3
L	21	LEU	-	expression tag	UNP Q6G4Y3
L	186	GLU	-	expression tag	UNP Q6G4Y3
L	187	GLY	-	expression tag	UNP Q6G4Y3
L	188	HIS	-	expression tag	UNP Q6G4Y3
L	189	HIS	-	expression tag	UNP Q6G4Y3
L	190	HIS	-	expression tag	UNP Q6G4Y3
L	191	HIS	-	expression tag	UNP Q6G4Y3
L	192	HIS	-	expression tag	UNP Q6G4Y3

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Chain	Residue	Modelled	Actual	Comment	Reference
L	193	HIS	-	expression tag	UNP Q6G4Y3

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		
2	H	1	Total	C	O	0	0
			6	3	3		
2	H	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	I	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	J	1	Total	C	O	0	0
			6	3	3		
2	K	1	Total	C	O	0	0
			6	3	3		
2	L	1	Total	C	O	0	0
			6	3	3		
2	L	1	Total	C	O	0	0
			6	3	3		

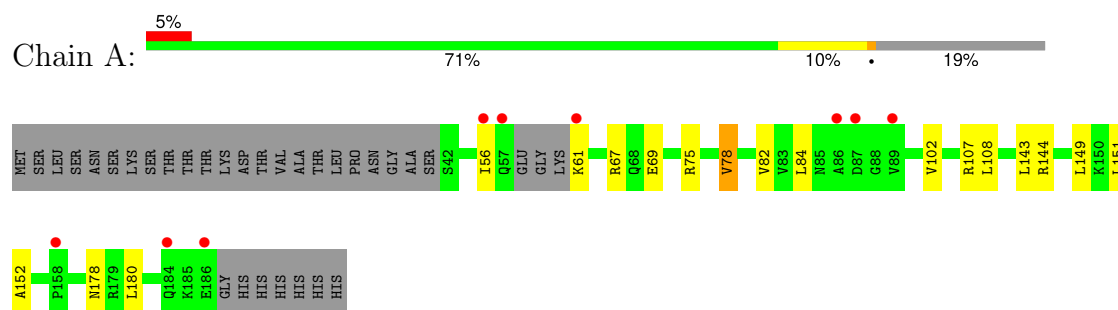
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	31	Total	O	0	0
			31	31		
3	B	46	Total	O	0	0
			46	46		
3	C	44	Total	O	0	0
			44	44		
3	D	35	Total	O	0	0
			35	35		
3	E	28	Total	O	0	0
			28	28		
3	F	44	Total	O	0	0
			44	44		
3	G	39	Total	O	0	0
			39	39		
3	H	25	Total	O	0	0
			25	25		
3	I	35	Total	O	0	0
			35	35		
3	J	32	Total	O	0	0
			32	32		
3	K	32	Total	O	0	0
			32	32		
3	L	25	Total	O	0	0
			25	25		

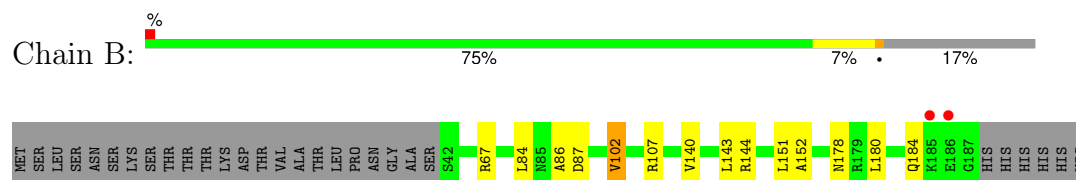
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

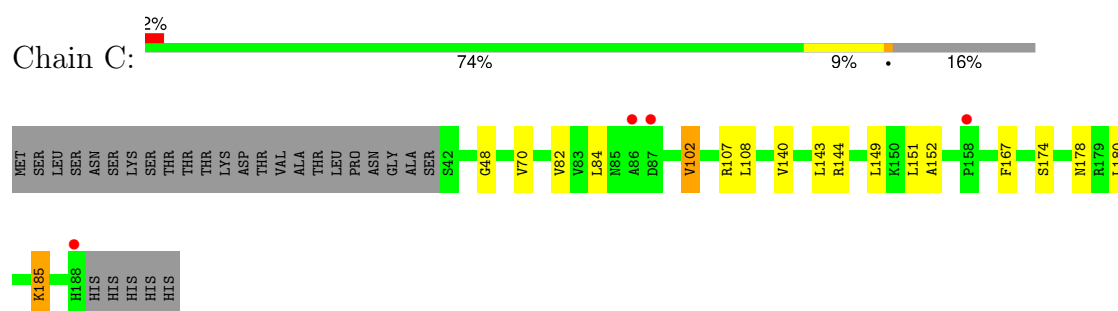
#### • Molecule 1: Invasion-associated protein B



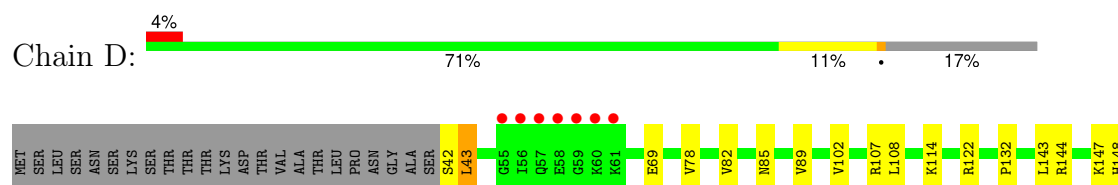
#### • Molecule 1: Invasion-associated protein B

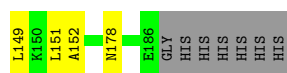


#### • Molecule 1: Invasion-associated protein B

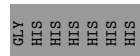
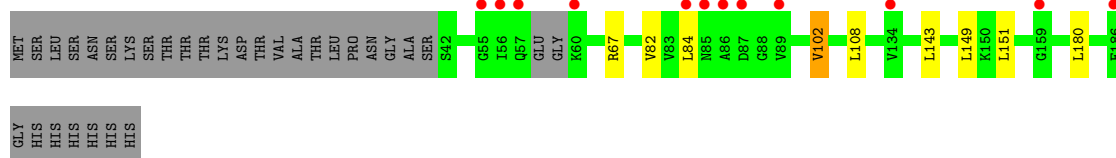
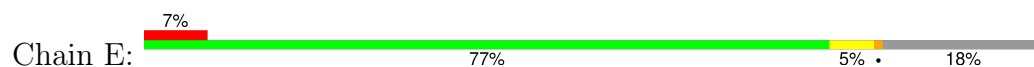


#### • Molecule 1: Invasion-associated protein B

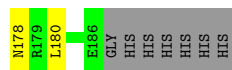
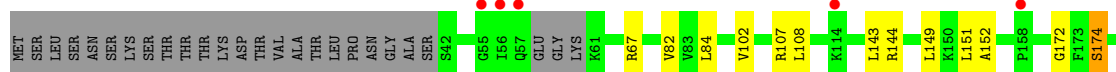




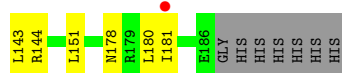
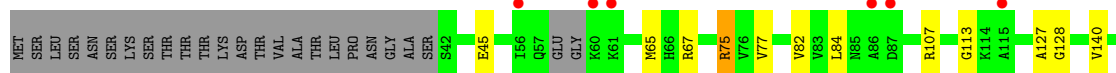
- Molecule 1: Invasion-associated protein B



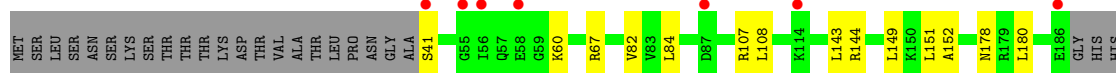
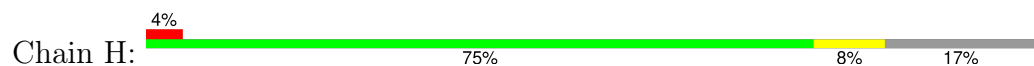
- Molecule 1: Invasion-associated protein B



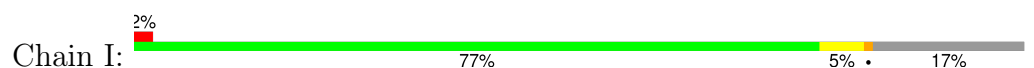
- Molecule 1: Invasion-associated protein B



- Molecule 1: Invasion-associated protein B



- Molecule 1: Invasion-associated protein B

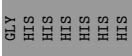
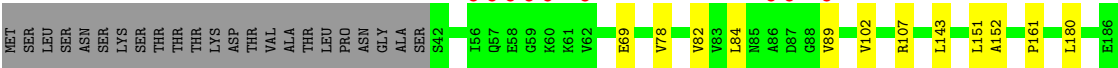
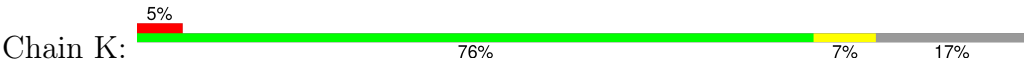




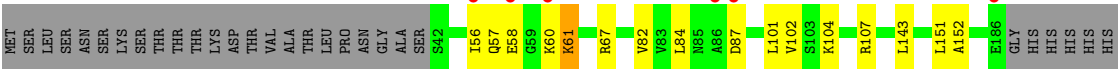
● Molecule 1: Invasion-associated protein B



● Molecule 1: Invasion-associated protein B



● Molecule 1: Invasion-associated protein B



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.71Å 139.95Å 179.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.35 36.76 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.3 (20.00-2.35) 99.3 (36.76-2.35)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.42 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.235 , 0.265 0.232 , 0.263	Depositor DCC
$R_{free}$ test set	3005 reflections (3.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	50.5	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 39.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13760	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.40	0/1090	0.61	0/1480
1	B	0.39	0/1136	0.61	0/1541
1	C	0.43	0/1134	0.63	0/1539
1	D	0.37	0/1132	0.61	0/1536
1	E	0.38	0/1110	0.59	0/1505
1	F	0.40	0/1112	0.62	0/1508
1	G	0.42	0/1116	0.63	0/1514
1	H	0.37	0/1138	0.61	0/1544
1	I	0.40	0/1131	0.60	0/1535
1	J	0.45	0/1130	0.61	0/1533
1	K	0.42	0/1135	0.61	0/1539
1	L	0.40	0/1130	0.61	0/1533
All	All	0.40	0/13494	0.61	0/18307

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	J	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	J	59	GLY	Peptide

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1074	0	1110	5	0
1	B	1113	0	1155	5	0
1	C	1113	0	1148	12	0
1	D	1109	0	1152	8	0
1	E	1091	0	1136	2	0
1	F	1090	0	1136	8	0
1	G	1097	0	1137	11	0
1	H	1115	0	1157	3	0
1	I	1108	0	1155	3	0
1	J	1107	0	1151	7	0
1	K	1112	0	1159	3	0
1	L	1107	0	1151	7	0
2	B	12	0	16	0	0
2	C	24	0	31	2	0
2	D	6	0	8	0	0
2	F	6	0	8	5	0
2	G	6	0	8	4	0
2	H	12	0	16	0	0
2	I	18	0	24	1	0
2	J	6	0	8	0	0
2	K	6	0	8	0	0
2	L	12	0	16	0	0
3	A	31	0	0	0	0
3	B	46	0	0	0	0
3	C	44	0	0	0	0
3	D	35	0	0	0	0
3	E	28	0	0	0	0
3	F	44	0	0	0	0
3	G	39	0	0	0	0
3	H	25	0	0	0	0
3	I	35	0	0	0	0
3	J	32	0	0	0	0
3	K	32	0	0	0	0
3	L	25	0	0	0	0
All	All	13760	0	13890	65	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 2.

All (65) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:185:LYS:HE2	1:J:60:LYS:HE3	1.29	1.09
1:L:61:LYS:O	1:L:61:LYS:HG2	1.66	0.95
1:C:185:LYS:CE	1:J:60:LYS:HE3	2.03	0.88
1:F:174:SER:H	2:F:1:GOL:H31	1.46	0.80
1:L:61:LYS:O	1:L:61:LYS:CG	2.30	0.80
1:F:174:SER:HB3	2:F:1:GOL:H32	1.63	0.79
1:C:185:LYS:HE2	1:J:60:LYS:CE	2.16	0.72
1:C:144:ARG:NH2	1:C:178:ASN:OD1	2.29	0.64
1:B:144:ARG:NH2	1:B:178:ASN:OD1	2.31	0.64
1:F:172:GLY:HA2	2:F:1:GOL:H12	1.80	0.63
1:F:172:GLY:CA	2:F:1:GOL:H12	2.28	0.62
1:F:144[B]:ARG:NH1	1:F:178:ASN:OD1	2.33	0.61
1:B:86:ALA:HA	1:B:184:GLN:HE22	1.67	0.58
1:I:122[A]:ARG:HB2	2:I:194:GOL:H2	1.87	0.55
1:C:167:PHE:CZ	1:G:45:GLU:HG2	2.41	0.55
1:F:174:SER:HB3	2:F:1:GOL:C3	2.35	0.55
1:A:144:ARG:NH2	1:A:178:ASN:OD1	2.40	0.54
1:G:181:ILE:HG21	1:L:60:LYS:HG3	1.91	0.53
1:G:77:VAL:O	2:G:1:GOL:H12	2.08	0.53
1:F:107:ARG:HB3	1:F:152:ALA:HB3	1.92	0.52
1:A:107:ARG:HB3	1:A:152:ALA:HB3	1.91	0.51
1:L:57:GLN:O	1:L:58:GLU:C	2.48	0.51
1:C:107:ARG:HB3	1:C:152:ALA:HB3	1.92	0.51
1:H:144:ARG:NH2	1:H:178:ASN:OD1	2.44	0.51
1:C:70:VAL:HG22	2:C:1:GOL:H32	1.93	0.50
1:C:102:VAL:HG13	1:E:102:VAL:HG13	1.93	0.50
1:A:69:GLU:HB2	1:A:78:VAL:HG13	1.93	0.50
1:C:140:VAL:O	1:C:144:ARG:HG3	2.12	0.50
1:C:167:PHE:HZ	1:G:45:GLU:HG2	1.78	0.48
1:L:56:ILE:O	1:L:56:ILE:HG13	2.13	0.48
1:D:147:LYS:O	1:D:148:HIS:HD2	1.96	0.48
1:B:107:ARG:HB3	1:B:152:ALA:HB3	1.97	0.47
1:F:108:LEU:HD23	1:F:149:LEU:HD21	1.97	0.47
1:G:144:ARG:NH2	1:G:178:ASN:OD1	2.47	0.46
1:I:107:ARG:HB3	1:I:152:ALA:HB3	1.98	0.46
1:D:85:ASN:HD22	1:D:89:VAL:HB	1.80	0.46
1:K:107:ARG:HB3	1:K:152:ALA:HB3	1.96	0.46
1:D:122[B]:ARG:HD2	1:D:132:PRO:HG3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:75:ARG:NH1	2:G:1:GOL:O2	2.50	0.45
1:L:101:LEU:HD23	1:L:104:LYS:HG3	1.97	0.44
1:H:108:LEU:HD23	1:H:149:LEU:HD21	1.99	0.44
1:D:144:ARG:NH2	1:D:178:ASN:OD1	2.50	0.44
1:J:107[B]:ARG:HG3	1:J:118:GLU:HG3	1.99	0.44
1:C:48:GLY:N	2:C:195:GOL:H2	2.33	0.44
1:B:140:VAL:O	1:B:144:ARG:HG3	2.17	0.43
1:E:108:LEU:HD23	1:E:149:LEU:HD21	1.99	0.43
1:D:107:ARG:HB3	1:D:152:ALA:HB3	2.00	0.43
1:B:102:VAL:HB	1:J:122:ARG:O	2.19	0.43
1:H:107:ARG:HB3	1:H:152:ALA:HB3	2.01	0.43
1:K:69:GLU:HB2	1:K:78:VAL:HG13	2.00	0.42
1:C:108:LEU:HD23	1:C:149:LEU:HD21	2.01	0.42
1:G:140:VAL:O	1:G:144:ARG:HG3	2.20	0.42
1:D:43:LEU:HD12	1:D:43:LEU:HA	1.79	0.42
1:A:56:ILE:HA	1:A:61:LYS:HA	2.02	0.41
1:G:65:MET:HB2	1:G:180:LEU:HD11	2.02	0.41
1:I:147:LYS:O	1:I:148:HIS:HD2	2.03	0.41
1:J:108:LEU:HD23	1:J:149:LEU:HD21	2.02	0.41
1:D:69:GLU:HB2	1:D:78:VAL:HG13	2.03	0.41
1:G:113:GLY:HA2	1:K:161:PRO:HA	2.02	0.41
1:G:127:ALA:O	2:G:1:GOL:H31	2.20	0.41
1:L:107:ARG:HB3	1:L:152:ALA:HB3	2.02	0.41
1:D:108:LEU:HD23	1:D:149:LEU:HD21	2.02	0.41
1:A:108:LEU:HD23	1:A:149:LEU:HD21	2.02	0.40
1:G:128:GLY:HA3	2:G:1:GOL:H31	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	138/175 (79%)	137 (99%)	1 (1%)	0	100	100
1	B	146/175 (83%)	143 (98%)	3 (2%)	0	100	100
1	C	146/175 (83%)	143 (98%)	3 (2%)	0	100	100
1	D	145/175 (83%)	142 (98%)	3 (2%)	0	100	100
1	E	140/175 (80%)	139 (99%)	1 (1%)	0	100	100
1	F	140/175 (80%)	140 (100%)	0	0	100	100
1	G	141/175 (81%)	140 (99%)	1 (1%)	0	100	100
1	H	146/175 (83%)	144 (99%)	2 (1%)	0	100	100
1	I	145/175 (83%)	145 (100%)	0	0	100	100
1	J	145/175 (83%)	142 (98%)	2 (1%)	1 (1%)	22	23
1	K	145/175 (83%)	142 (98%)	3 (2%)	0	100	100
1	L	145/175 (83%)	141 (97%)	4 (3%)	0	100	100
All	All	1722/2100 (82%)	1698 (99%)	23 (1%)	1 (0%)	51	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	J	59	GLY

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	119/147 (81%)	110 (92%)	9 (8%)	13	13
1	B	123/147 (84%)	116 (94%)	7 (6%)	20	22
1	C	123/147 (84%)	115 (94%)	8 (6%)	17	18
1	D	123/147 (84%)	116 (94%)	7 (6%)	20	22
1	E	121/147 (82%)	114 (94%)	7 (6%)	20	22
1	F	121/147 (82%)	113 (93%)	8 (7%)	16	17
1	G	122/147 (83%)	115 (94%)	7 (6%)	20	22

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	124/147 (84%)	116 (94%)	8 (6%)	17	18
1	I	123/147 (84%)	116 (94%)	7 (6%)	20	22
1	J	123/147 (84%)	116 (94%)	7 (6%)	20	22
1	K	123/147 (84%)	116 (94%)	7 (6%)	20	22
1	L	123/147 (84%)	115 (94%)	8 (6%)	17	18
All	All	1468/1764 (83%)	1378 (94%)	90 (6%)	18	20

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	ARG
1	A	75	ARG
1	A	78	VAL
1	A	82	VAL
1	A	84	LEU
1	A	102	VAL
1	A	143	LEU
1	A	151	LEU
1	A	180	LEU
1	B	67	ARG
1	B	84	LEU
1	B	87	ASP
1	B	102	VAL
1	B	143	LEU
1	B	151	LEU
1	B	180	LEU
1	C	82	VAL
1	C	84	LEU
1	C	102	VAL
1	C	143	LEU
1	C	151	LEU
1	C	174	SER
1	C	180	LEU
1	C	185	LYS
1	D	42	SER
1	D	43	LEU
1	D	82	VAL
1	D	102	VAL
1	D	114	LYS
1	D	143	LEU

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Mol	Chain	Res	Type
1	D	151	LEU
1	E	67	ARG
1	E	82	VAL
1	E	84	LEU
1	E	102	VAL
1	E	143	LEU
1	E	151	LEU
1	E	180	LEU
1	F	67	ARG
1	F	82	VAL
1	F	84	LEU
1	F	102	VAL
1	F	143	LEU
1	F	151	LEU
1	F	174	SER
1	F	180	LEU
1	G	67	ARG
1	G	75	ARG
1	G	82	VAL
1	G	84	LEU
1	G	107	ARG
1	G	143	LEU
1	G	151	LEU
1	H	41	SER
1	H	60	LYS
1	H	67	ARG
1	H	82	VAL
1	H	84	LEU
1	H	143	LEU
1	H	151	LEU
1	H	180	LEU
1	I	67	ARG
1	I	84	LEU
1	I	102	VAL
1	I	122[A]	ARG
1	I	122[B]	ARG
1	I	143	LEU
1	I	151	LEU
1	J	60	LYS
1	J	82	VAL
1	J	84	LEU
1	J	143	LEU

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Mol	Chain	Res	Type
1	J	150	LYS
1	J	151	LEU
1	J	180	LEU
1	K	82	VAL
1	K	84	LEU
1	K	89	VAL
1	K	102	VAL
1	K	143	LEU
1	K	151	LEU
1	K	180	LEU
1	L	61	LYS
1	L	67	ARG
1	L	82	VAL
1	L	84	LEU
1	L	87	ASP
1	L	102	VAL
1	L	143	LEU
1	L	151	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	B	85	ASN
1	B	148	HIS
1	B	184	GLN
1	C	68	GLN
1	C	85	ASN
1	C	148	HIS
1	D	53	ASN
1	D	85	ASN
1	D	148	HIS
1	E	85	ASN
1	E	148	HIS
1	F	148	HIS
1	G	148	HIS
1	H	57	GLN
1	H	148	HIS
1	I	148	HIS
1	J	148	HIS
1	J	184	GLN
1	K	148	HIS

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Mol	Chain	Res	Type
1	L	148	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

18 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	GOL	D	1	-	5,5,5	0.39	0	5,5,5	0.39	0
2	GOL	C	196	-	5,5,5	0.41	0	5,5,5	0.35	0
2	GOL	I	194	-	5,5,5	0.37	0	5,5,5	0.23	0
2	GOL	L	1	-	5,5,5	0.37	0	5,5,5	0.28	0
2	GOL	I	1	-	5,5,5	0.34	0	5,5,5	0.51	0
2	GOL	J	1	-	5,5,5	0.76	0	5,5,5	0.55	0
2	GOL	G	1	-	5,5,5	0.77	0	5,5,5	1.00	0
2	GOL	F	1	-	5,5,5	0.74	0	5,5,5	0.42	0
2	GOL	H	1	-	5,5,5	0.39	0	5,5,5	0.27	0
2	GOL	L	194	-	5,5,5	0.38	0	5,5,5	0.25	0
2	GOL	H	194	-	5,5,5	0.37	0	5,5,5	0.30	0
2	GOL	K	1	-	5,5,5	0.36	0	5,5,5	0.34	0
2	GOL	B	1	-	5,5,5	0.38	0	5,5,5	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	GOL	B	194	-	5,5,5	0.36	0	5,5,5	0.32	0
2	GOL	C	1	-	5,5,5	0.38	0	5,5,5	0.37	0
2	GOL	I	195	-	5,5,5	0.37	0	5,5,5	0.35	0
2	GOL	C	194	-	5,5,5	0.37	0	5,5,5	0.59	0
2	GOL	C	195	-	5,5,5	1.08	1 (20%)	5,5,5	0.72	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	D	1	-	-	0/4/4/4	-
2	GOL	C	196	-	-	2/4/4/4	-
2	GOL	I	194	-	-	2/4/4/4	-
2	GOL	L	1	-	-	4/4/4/4	-
2	GOL	I	1	-	-	2/4/4/4	-
2	GOL	J	1	-	-	0/4/4/4	-
2	GOL	G	1	-	-	4/4/4/4	-
2	GOL	F	1	-	-	2/4/4/4	-
2	GOL	H	1	-	-	2/4/4/4	-
2	GOL	L	194	-	-	4/4/4/4	-
2	GOL	H	194	-	-	2/4/4/4	-
2	GOL	K	1	-	-	2/4/4/4	-
2	GOL	B	1	-	-	4/4/4/4	-
2	GOL	B	194	-	-	4/4/4/4	-
2	GOL	C	1	-	-	2/4/4/4	-
2	GOL	I	195	-	-	4/4/4/4	-
2	GOL	C	194	-	-	3/4/4/4	-
2	GOL	C	195	-	-	4/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	195	GOL	O2-C2	-2.26	1.36	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	GOL	C1-C2-C3-O3
2	B	194	GOL	O1-C1-C2-C3
2	C	1	GOL	O1-C1-C2-O2
2	C	1	GOL	O1-C1-C2-C3
2	C	195	GOL	O1-C1-C2-C3
2	C	196	GOL	C1-C2-C3-O3
2	C	196	GOL	O2-C2-C3-O3
2	G	1	GOL	O1-C1-C2-O2
2	G	1	GOL	C1-C2-C3-O3
2	H	1	GOL	C1-C2-C3-O3
2	H	194	GOL	O1-C1-C2-O2
2	H	194	GOL	O1-C1-C2-C3
2	I	1	GOL	O1-C1-C2-C3
2	I	194	GOL	O1-C1-C2-C3
2	I	195	GOL	O1-C1-C2-C3
2	K	1	GOL	O1-C1-C2-C3
2	L	1	GOL	O1-C1-C2-O2
2	L	1	GOL	O1-C1-C2-C3
2	L	1	GOL	C1-C2-C3-O3
2	L	194	GOL	O1-C1-C2-C3
2	L	194	GOL	C1-C2-C3-O3
2	B	1	GOL	O2-C2-C3-O3
2	B	194	GOL	O1-C1-C2-O2
2	I	1	GOL	O1-C1-C2-O2
2	L	194	GOL	O1-C1-C2-O2
2	B	1	GOL	O1-C1-C2-C3
2	B	194	GOL	C1-C2-C3-O3
2	C	194	GOL	C1-C2-C3-O3
2	C	195	GOL	C1-C2-C3-O3
2	G	1	GOL	O1-C1-C2-C3
2	I	195	GOL	C1-C2-C3-O3
2	B	194	GOL	O2-C2-C3-O3
2	C	195	GOL	O1-C1-C2-O2
2	C	195	GOL	O2-C2-C3-O3
2	F	1	GOL	O2-C2-C3-O3
2	G	1	GOL	O2-C2-C3-O3
2	H	1	GOL	O2-C2-C3-O3
2	I	194	GOL	O1-C1-C2-O2
2	I	195	GOL	O1-C1-C2-O2
2	K	1	GOL	O1-C1-C2-O2
2	L	1	GOL	O2-C2-C3-O3
2	I	195	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
2	L	194	GOL	O2-C2-C3-O3
2	B	1	GOL	O1-C1-C2-O2
2	C	194	GOL	O2-C2-C3-O3
2	F	1	GOL	C1-C2-C3-O3
2	C	194	GOL	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	194	GOL	1	0
2	G	1	GOL	4	0
2	F	1	GOL	5	0
2	C	1	GOL	1	0
2	C	195	GOL	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å <sup>2</sup> )	Q < 0.9
1	A	142/175 (81%)	0.51	9 (6%)	20 29	32, 46, 80, 108	0
1	B	146/175 (83%)	0.26	2 (1%)	75 83	32, 45, 73, 87	0
1	C	147/175 (84%)	0.40	4 (2%)	54 64	31, 45, 73, 108	0
1	D	145/175 (82%)	0.45	7 (4%)	30 43	31, 46, 83, 111	0
1	E	143/175 (81%)	0.41	12 (8%)	11 16	32, 45, 80, 116	0
1	F	142/175 (81%)	0.36	5 (3%)	44 56	31, 45, 71, 104	0
1	G	143/175 (81%)	0.38	7 (4%)	29 42	32, 46, 77, 98	0
1	H	146/175 (83%)	0.18	7 (4%)	30 43	32, 47, 81, 113	0
1	I	145/175 (82%)	0.28	3 (2%)	63 74	32, 46, 78, 86	0
1	J	145/175 (82%)	0.33	6 (4%)	37 49	31, 46, 77, 96	0
1	K	145/175 (82%)	0.31	9 (6%)	20 29	32, 46, 91, 119	0
1	L	145/175 (82%)	0.28	6 (4%)	37 49	32, 46, 92, 128	0
All	All	1734/2100 (82%)	0.35	77 (4%)	34 46	31, 46, 81, 128	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	86	ALA	8.5
1	K	86	ALA	6.4
1	D	56	ILE	5.9
1	K	56	ILE	5.9
1	E	56	ILE	5.6
1	L	86	ALA	5.1
1	G	60	LYS	4.7
1	E	87	ASP	4.7
1	K	87	ASP	4.4
1	E	86	ALA	4.3
1	H	41	SER	4.3

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Mol	Chain	Res	Type	RSRZ
1	L	56	ILE	4.2
1	D	58	GLU	4.0
1	L	58	GLU	4.0
1	I	186	GLU	4.0
1	E	60	LYS	3.9
1	H	58	GLU	3.9
1	J	159	GLY	3.8
1	J	158	PRO	3.8
1	L	87	ASP	3.7
1	E	186	GLU	3.7
1	D	61	LYS	3.7
1	C	188	HIS	3.6
1	K	58	GLU	3.6
1	D	59	GLY	3.3
1	A	56	ILE	3.3
1	K	60	LYS	3.3
1	G	86	ALA	3.2
1	H	56	ILE	3.1
1	A	87	ASP	3.1
1	I	86	ALA	3.1
1	F	57	GLN	3.1
1	G	56	ILE	3.0
1	F	56	ILE	3.0
1	H	87	ASP	3.0
1	K	62	VAL	2.9
1	H	55	GLY	2.9
1	A	86	ALA	2.8
1	A	158	PRO	2.8
1	D	55	GLY	2.7
1	I	58	GLU	2.7
1	G	87	ASP	2.6
1	J	186	GLU	2.6
1	A	186	GLU	2.6
1	L	186	GLU	2.6
1	A	184	GLN	2.5
1	D	60	LYS	2.5
1	E	84	LEU	2.5
1	E	89	VAL	2.5
1	E	134	VAL	2.5
1	K	57	GLN	2.4
1	F	158	PRO	2.4
1	B	186	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	J	85	ASN	2.4
1	C	87	ASP	2.4
1	G	61	LYS	2.4
1	K	89	VAL	2.4
1	E	57	GLN	2.3
1	F	55	GLY	2.3
1	E	85	ASN	2.3
1	H	186	GLU	2.3
1	H	114	LYS	2.3
1	J	60	LYS	2.3
1	F	114	LYS	2.2
1	E	159	GLY	2.2
1	L	60	LYS	2.2
1	C	158	PRO	2.2
1	A	89	VAL	2.2
1	B	185	LYS	2.2
1	A	61	LYS	2.2
1	D	57	GLN	2.2
1	A	57	GLN	2.2
1	K	59	GLY	2.2
1	G	115	ALA	2.1
1	C	86	ALA	2.1
1	E	55	GLY	2.1
1	G	181	ILE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GOL	L	194	6/6	0.70	0.26	81,88,89,92	0
2	GOL	D	1	6/6	0.72	0.41	54,69,71,78	0
2	GOL	B	194	6/6	0.72	0.31	73,75,79,84	0
2	GOL	I	1	6/6	0.76	0.30	65,70,73,81	0
2	GOL	B	1	6/6	0.81	0.31	75,84,86,87	0
2	GOL	H	194	6/6	0.81	0.23	66,73,78,78	0
2	GOL	C	194	6/6	0.82	0.28	46,53,60,63	0
2	GOL	J	1	6/6	0.83	0.33	55,70,75,75	0
2	GOL	C	1	6/6	0.83	0.33	45,52,59,69	0
2	GOL	C	195	6/6	0.84	0.42	34,61,63,66	0
2	GOL	I	194	6/6	0.84	0.23	73,81,84,88	0
2	GOL	L	1	6/6	0.85	0.40	70,77,78,82	0
2	GOL	H	1	6/6	0.85	0.40	61,70,75,78	0
2	GOL	F	1	6/6	0.86	0.33	49,53,60,68	0
2	GOL	C	196	6/6	0.86	0.26	68,77,85,87	0
2	GOL	I	195	6/6	0.88	0.31	55,60,63,65	0
2	GOL	K	1	6/6	0.89	0.25	72,75,76,79	0
2	GOL	G	1	6/6	0.91	0.30	64,77,81,83	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.