



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2021 – 10:10 AM JST

PDB ID : 7BPK  
Title : Zika virus envelope protein mutant bound to mAb  
Authors : Dai, L.; Qi, J.; Gao, G.F.  
Deposited on : 2020-03-23  
Resolution : 3.10 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.

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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
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
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)	:	2.23.2

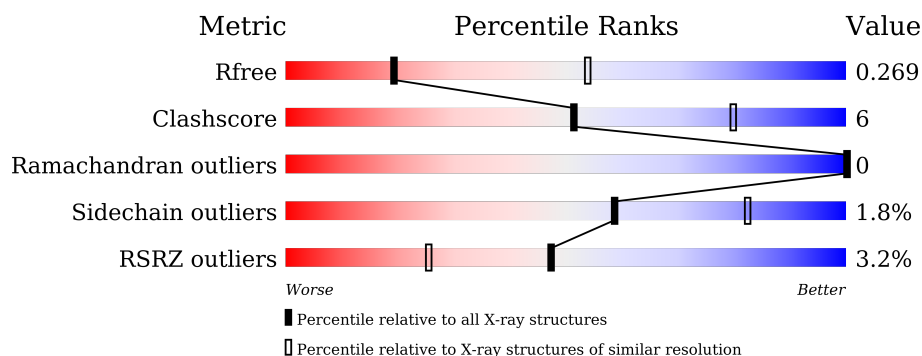
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	416	<div> <div>6%</div> <div> <div></div> <div>80%</div> <div>17%</div> <div></div> </div> <div></div> </div>
1	B	416	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div></div> </div> <div></div> </div>
2	C	127	<div> <div>7%</div> <div> <div></div> <div>80%</div> <div>20%</div> <div></div> </div> <div></div> </div>
2	H	127	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>24%</div> <div></div> </div> <div></div> </div>
3	D	108	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>12%</div> <div></div> </div> <div></div> </div>
3	L	108	<div> <div></div> <div> <div></div> <div>88%</div> <div>12%</div> <div></div> </div> <div></div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9754 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 Envelope protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			3094	1931	540	596	27			
1	B	404	Total	C	N	O	S	0	0	0
			3094	1931	540	596	27			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP A0A142I5B9
A	98	ASN	ASP	conflict	UNP A0A142I5B9
A	103	THR	ASN	conflict	UNP A0A142I5B9
A	106	LEU	GLY	conflict	UNP A0A142I5B9
A	107	GLU	LEU	conflict	UNP A0A142I5B9
A	108	TRP	PHE	conflict	UNP A0A142I5B9
A	410	HIS	-	expression tag	UNP A0A142I5B9
A	411	HIS	-	expression tag	UNP A0A142I5B9
A	412	HIS	-	expression tag	UNP A0A142I5B9
A	413	HIS	-	expression tag	UNP A0A142I5B9
A	414	HIS	-	expression tag	UNP A0A142I5B9
A	415	HIS	-	expression tag	UNP A0A142I5B9
B	0	MET	-	expression tag	UNP A0A142I5B9
B	98	ASN	ASP	conflict	UNP A0A142I5B9
B	103	THR	ASN	conflict	UNP A0A142I5B9
B	106	LEU	GLY	conflict	UNP A0A142I5B9
B	107	GLU	LEU	conflict	UNP A0A142I5B9
B	108	TRP	PHE	conflict	UNP A0A142I5B9
B	410	HIS	-	expression tag	UNP A0A142I5B9
B	411	HIS	-	expression tag	UNP A0A142I5B9
B	412	HIS	-	expression tag	UNP A0A142I5B9
B	413	HIS	-	expression tag	UNP A0A142I5B9
B	414	HIS	-	expression tag	UNP A0A142I5B9
B	415	HIS	-	expression tag	UNP A0A142I5B9

- Molecule 2 is a protein called Z3L1 Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	127	Total	C	N	O	S	0	0	0
			991	627	170	190	4			
2	H	127	Total	C	N	O	S	0	0	0
			991	627	170	190	4			

- Molecule 3 is a protein called IG c307\_light\_IGLV1-51\_IGLJ2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	108	Total	C	N	O	S	0	0	0
			792	495	132	163	2			
3	L	108	Total	C	N	O	S	0	0	0
			792	495	132	163	2			

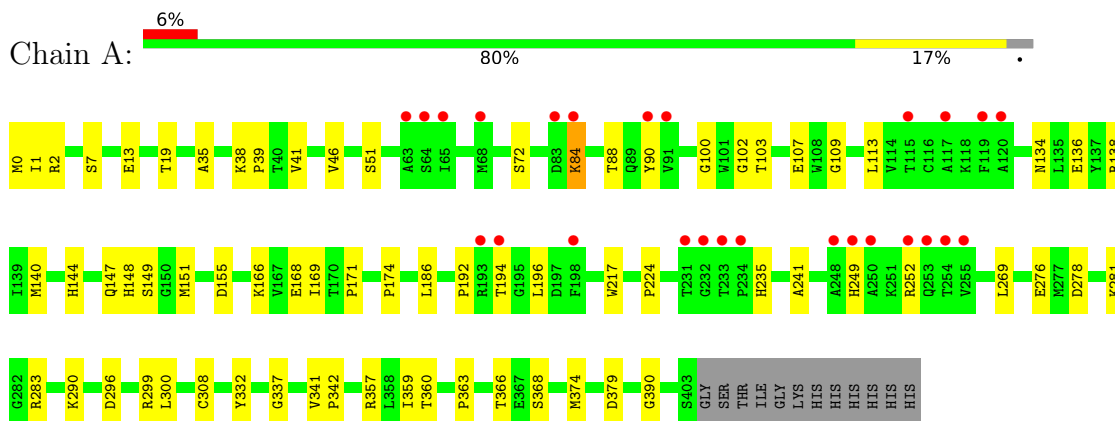
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	99	TRP	LEU	conflict	UNP A0A5C2GK82
L	99	TRP	LEU	conflict	UNP A0A5C2GK82

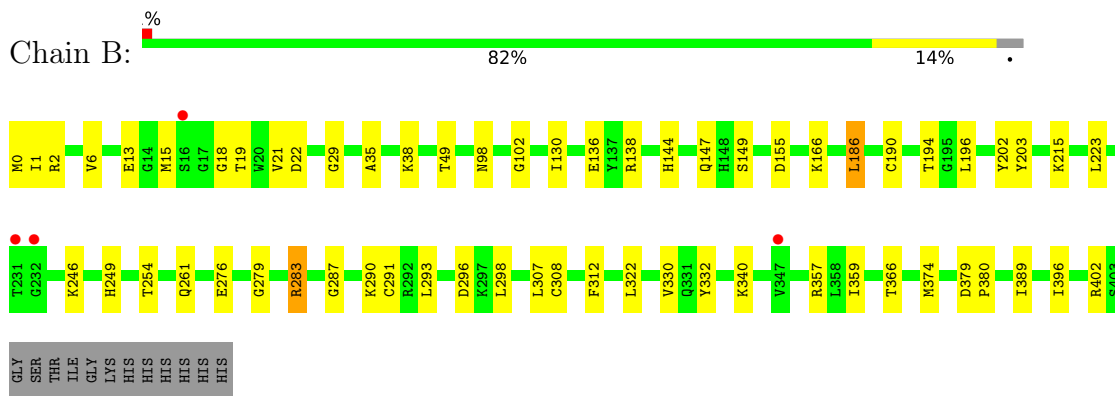
### 3 Residue-property plots [i](#)

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.

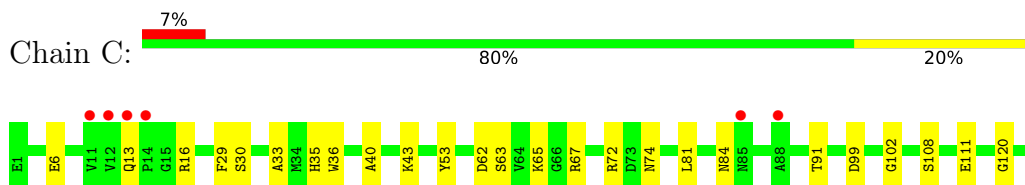
- Molecule 1: Envelope protein



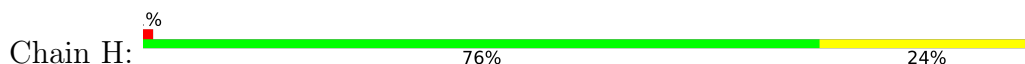
- Molecule 1: Envelope protein

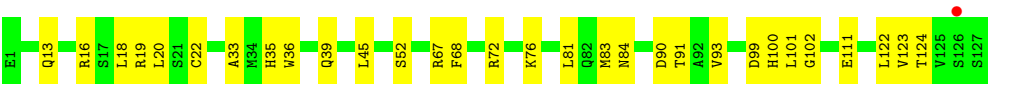


- Molecule 2: Z3L1 Heavy chain

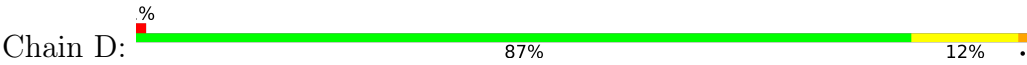


- Molecule 2: Z3L1 Heavy chain

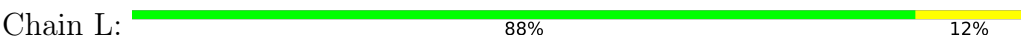




● Molecule 3: IG c307\_light\_IGLV1-51\_IGLJ2



● Molecule 3: IG c307\_light\_IGLV1-51\_IGLJ2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.46Å 103.79Å 205.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.45 – 3.10 36.45 – 3.10	Depositor EDS
% Data completeness (in resolution range)	95.4 (36.45-3.10) 95.4 (36.45-3.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.25 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, $R_{free}$	0.232 , 0.268 0.233 , 0.269	Depositor DCC
$R_{free}$ test set	1378 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	69.6	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 37.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	9754	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% 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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.24	0/3160	0.45	0/4283
1	B	0.24	0/3160	0.45	0/4283
2	C	0.24	0/1017	0.44	0/1381
2	H	0.24	0/1017	0.46	0/1381
3	D	0.24	0/811	0.45	0/1106
3	L	0.25	0/811	0.44	0/1106
All	All	0.24	0/9976	0.45	0/13540

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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	3094	0	3024	42	0
1	B	3094	0	3024	34	0
2	C	991	0	942	15	0
2	H	991	0	942	16	0
3	D	792	0	765	8	0
3	L	792	0	765	5	0
All	All	9754	0	9462	108	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 6.



All (108) 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:A:138:ARG:NH2	2:C:111:GLU:OE2	2.23	0.72
1:A:102:GLY:HA3	1:B:1:ILE:HG21	1.76	0.67
1:A:147:GLN:OE1	1:A:366:THR:N	2.22	0.65
1:A:51:SER:HB3	1:A:281:LYS:HG2	1.77	0.65
2:C:29:PHE:O	2:C:72:ARG:NH2	2.31	0.64
3:L:62:ARG:NH1	3:L:83:ASP:OD2	2.31	0.63
1:B:194:THR:HG22	1:B:287:GLY:HA2	1.81	0.62
3:L:32:ASN:ND2	3:L:92:TRP:O	2.26	0.61
1:A:35:ALA:HB3	1:A:38:LYS:HB2	1.81	0.61
1:B:138:ARG:NH2	2:H:111:GLU:OE2	2.33	0.61
2:H:91:THR:HG23	2:H:124:THR:HA	1.85	0.59
2:H:13:GLN:H	2:H:16:ARG:HH21	1.51	0.58
1:B:19:THR:HG22	1:B:296:ASP:HB3	1.86	0.58
3:L:40:LEU:HD23	3:L:85:ALA:HB2	1.86	0.57
1:A:249:HIS:HB2	1:A:252:ARG:HG2	1.86	0.57
1:A:0:MET:HG3	1:A:149:SER:HB2	1.86	0.57
2:H:67:ARG:NH2	2:H:90:ASP:OD2	2.37	0.57
1:A:194:THR:HG21	1:A:290:LYS:HB2	1.85	0.56
1:A:1:ILE:HG21	1:B:102:GLY:HA3	1.87	0.56
2:C:62:ASP:HA	2:C:65:LYS:HG3	1.87	0.56
3:D:32:ASN:ND2	3:D:92:TRP:O	2.33	0.56
1:B:276:GLU:HB3	1:B:283:ARG:HB3	1.86	0.56
3:D:40:LEU:HD23	3:D:85:ALA:HB2	1.86	0.56
2:H:52:SER:O	2:H:72:ARG:NH1	2.37	0.55
1:A:109:GLY:HA2	1:B:322:LEU:HD21	1.88	0.55
1:A:147:GLN:HA	1:A:374:MET:HG3	1.88	0.55
1:A:359:ILE:HD11	1:A:379:ASP:HB2	1.89	0.55
3:D:38:GLN:HB2	3:D:48:LEU:HD12	1.89	0.54
1:B:283:ARG:NH1	2:H:111:GLU:OE1	2.40	0.54
1:B:0:MET:HG3	1:B:144:HIS:CE1	2.43	0.54
1:A:224:PRO:HD3	1:A:241:ALA:HB3	1.90	0.53
1:B:147:GLN:OE1	1:B:366:THR:N	2.37	0.53
1:B:49:THR:HG23	1:B:136:GLU:HB3	1.89	0.53
1:B:35:ALA:HB3	1:B:38:LYS:HB2	1.91	0.53
2:H:35:HIS:CD2	2:H:99:ASP:HB2	2.44	0.53
1:A:276:GLU:HB2	1:A:283:ARG:HB3	1.90	0.52
3:L:17:LYS:HD3	3:L:77:THR:HG22	1.90	0.52
1:A:51:SER:O	1:A:134:ASN:ND2	2.37	0.52
1:B:223:LEU:HD21	1:B:261:GLN:HG3	1.91	0.51
1:A:41:VAL:HG21	1:A:186:LEU:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:SER:HB3	1:A:113:LEU:HD13	1.92	0.51
1:B:246:LYS:HB2	1:B:254:THR:OG1	2.11	0.51
1:B:279:GLY:HA3	2:H:101:LEU:HA	1.92	0.50
1:A:341:VAL:HB	1:A:363:PRO:HG2	1.94	0.50
2:H:36:TRP:NE1	2:H:81:LEU:HB2	2.27	0.50
1:B:21:VAL:HG22	1:B:293:LEU:HB2	1.94	0.50
1:A:7:SER:HB2	1:B:98:ASN:HD21	1.77	0.49
1:A:169:ILE:HG23	1:A:174:PRO:HA	1.94	0.49
2:C:91:THR:HG23	2:C:124:THR:HA	1.93	0.49
2:H:93:VAL:HG22	2:H:122:LEU:HD13	1.95	0.49
1:B:190:CYS:HA	1:B:291:CYS:HA	1.95	0.48
2:H:39:GLN:HB2	2:H:45:LEU:HD23	1.95	0.48
1:A:166:LYS:NZ	3:D:31:ASN:O	2.47	0.48
1:B:130:ILE:HD11	1:B:203:TYR:HB2	1.95	0.48
2:C:30:SER:HB3	2:C:74:ASN:HB3	1.96	0.48
1:A:100:GLY:N	1:A:103:THR:OG1	2.46	0.47
1:A:148:HIS:O	1:A:151:MET:HG2	2.14	0.47
2:C:35:HIS:CD2	2:C:99:ASP:HB2	2.49	0.47
3:L:93:ASP:OD2	3:L:96:LEU:HB2	2.14	0.47
1:B:147:GLN:O	1:B:147:GLN:HG2	2.15	0.47
1:B:6:VAL:HG11	1:B:29:GLY:HA2	1.96	0.46
1:B:2:ARG:HH22	1:B:155:ASP:CG	2.19	0.46
1:B:312:PHE:O	1:B:396:ILE:HD11	2.16	0.46
1:B:330:VAL:HG11	1:B:389:ILE:HD13	1.96	0.46
1:A:342:PRO:HG2	1:A:390:GLY:HA2	1.98	0.45
2:H:20:LEU:HD21	2:H:123:VAL:HG21	1.97	0.45
3:D:43:THR:HG22	3:D:44:ALA:H	1.82	0.45
1:A:84:LYS:HE3	1:A:90:TYR:CE2	2.51	0.45
3:D:89:CYS:O	3:D:102:GLY:N	2.49	0.45
1:B:307:LEU:HD23	1:B:340:LYS:HB3	1.98	0.45
2:H:99:ASP:OD1	2:H:100:HIS:N	2.50	0.45
3:D:62:ARG:NH1	3:D:83:ASP:OD2	2.50	0.45
1:B:15:MET:HB2	1:B:18:GLY:H	1.81	0.45
1:A:19:THR:HG22	1:A:296:ASP:HB3	1.98	0.45
1:B:308:CYS:HB3	1:B:332:TYR:CE1	2.52	0.44
2:C:33:ALA:H	2:C:102:GLY:HA2	1.80	0.44
1:A:144:HIS:HD2	1:A:360:THR:HG22	1.82	0.44
1:A:217:TRP:CH2	1:A:269:LEU:HD21	2.52	0.44
1:A:46:VAL:HG21	1:A:140:MET:HB2	1.99	0.43
2:C:63:SER:O	2:C:67:ARG:NH1	2.50	0.43
1:B:13:GLU:OE2	1:B:357:ARG:NH1	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:GLN:HA	1:B:374:MET:HG3	2.00	0.43
1:B:196:LEU:HD12	1:B:196:LEU:HA	1.83	0.43
1:A:278:ASP:OD2	2:C:108:SER:HB3	2.18	0.43
1:A:39:PRO:HD3	1:A:300:LEU:HD22	2.00	0.43
1:B:380:PRO:HD2	1:B:402:ARG:HH11	1.84	0.43
1:B:202:TYR:CE1	1:B:215:LYS:HE2	2.54	0.43
2:C:36:TRP:CG	2:C:81:LEU:HD22	2.54	0.43
2:C:36:TRP:NE1	2:C:81:LEU:HB2	2.35	0.42
1:A:136:GLU:OE2	1:A:168:GLU:HG2	2.20	0.42
1:A:171:PRO:HA	1:A:192:PRO:HG2	2.02	0.41
1:A:308:CYS:HB3	1:A:332:TYR:CZ	2.55	0.41
2:C:40:ALA:HB3	2:C:43:LYS:HB2	2.02	0.41
1:A:136:GLU:OE2	3:D:33:TYR:OH	2.39	0.41
1:A:196:LEU:HD12	1:A:196:LEU:HA	1.77	0.41
1:A:138:ARG:HG3	1:A:168:GLU:HG3	2.01	0.41
1:A:337:GLY:HA2	1:A:368:SER:HB3	2.02	0.41
1:A:13:GLU:OE1	1:A:357:ARG:NH1	2.48	0.41
1:B:359:ILE:HD11	1:B:379:ASP:HB2	2.03	0.41
2:H:33:ALA:N	2:H:102:GLY:HA2	2.36	0.41
1:A:283:ARG:NH1	2:C:111:GLU:OE1	2.54	0.40
1:B:186:LEU:HG	1:B:298:LEU:HD11	2.03	0.40
2:C:6:GLU:OE1	2:C:120:GLY:N	2.46	0.40
1:A:2:ARG:NH1	1:A:155:ASP:OD1	2.54	0.40
1:A:88:THR:HB	1:A:235:HIS:CE1	2.56	0.40
2:C:13:GLN:OE1	2:C:16:ARG:NH2	2.54	0.40
2:H:18:LEU:HD23	2:H:19:ARG:N	2.36	0.40
2:H:68:PHE:CE1	2:H:83:MET:HB3	2.57	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	402/416 (97%)	402 (100%)	0	0	100	100
1	B	402/416 (97%)	402 (100%)	0	0	100	100
2	C	125/127 (98%)	125 (100%)	0	0	100	100
2	H	125/127 (98%)	125 (100%)	0	0	100	100
3	D	106/108 (98%)	106 (100%)	0	0	100	100
3	L	106/108 (98%)	106 (100%)	0	0	100	100
All	All	1266/1302 (97%)	1266 (100%)	0	0	100	100

There are no Ramachandran outliers to report.

### 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	339/349 (97%)	336 (99%)	3 (1%)	78	91
1	B	339/349 (97%)	332 (98%)	7 (2%)	53	79
2	C	105/105 (100%)	103 (98%)	2 (2%)	57	81
2	H	105/105 (100%)	102 (97%)	3 (3%)	42	72
3	D	89/89 (100%)	88 (99%)	1 (1%)	73	89
3	L	89/89 (100%)	86 (97%)	3 (3%)	37	69
All	All	1066/1086 (98%)	1047 (98%)	19 (2%)	59	82

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	LYS
1	A	107	GLU
1	A	299	ARG
1	B	22	ASP
1	B	149	SER
1	B	166	LYS
1	B	186	LEU

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Mol	Chain	Res	Type
1	B	249	HIS
1	B	283	ARG
1	B	290	LYS
2	C	53	TYR
2	C	84	ASN
3	D	89	CYS
2	H	22	CYS
2	H	76	LYS
2	H	84	ASN
3	L	48	LEU
3	L	89	CYS
3	L	99	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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](#)

There are no ligands in this entry.

## 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	404/416 (97%)	0.23	26 (6%) 19 8	33, 70, 123, 187	0
1	B	404/416 (97%)	-0.02	4 (0%) 82 67	31, 65, 105, 165	0
2	C	127/127 (100%)	0.34	9 (7%) 16 6	50, 82, 121, 138	0
2	H	127/127 (100%)	-0.18	1 (0%) 86 72	29, 50, 76, 103	0
3	D	108/108 (100%)	0.05	1 (0%) 84 69	35, 67, 107, 115	0
3	L	108/108 (100%)	-0.14	0 100 100	24, 50, 85, 101	0
All	All	1278/1302 (98%)	0.07	41 (3%) 47 25	24, 66, 115, 187	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	232	GLY	5.4
2	C	127	SER	4.9
1	A	231	THR	4.6
1	A	232	GLY	4.4
1	A	233	THR	4.3
1	A	117	ALA	4.0
1	A	253	GLN	3.5
1	A	65	ILE	3.3
2	C	126	SER	3.2
1	A	64	SER	3.2
1	A	194	THR	3.1
1	A	250	ALA	3.0
1	A	63	ALA	2.9
1	A	120	ALA	2.9
1	A	68	MET	2.8
1	A	248	ALA	2.8
1	A	249	HIS	2.8
1	A	252	ARG	2.7
1	A	115	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	16	SER	2.6
1	A	198	PHE	2.6
2	C	11	VAL	2.5
2	C	13	GLN	2.4
2	C	14	PRO	2.4
1	A	234	PRO	2.3
2	C	125	VAL	2.3
1	A	90	TYR	2.3
1	A	91	VAL	2.3
1	B	231	THR	2.2
2	C	88	ALA	2.2
1	A	84	LYS	2.2
1	A	193	ARG	2.2
2	C	85	ASN	2.2
1	A	255	VAL	2.1
2	C	12	VAL	2.1
1	B	347	VAL	2.1
1	A	254	THR	2.1
3	D	106	LYS	2.1
1	A	119	PHE	2.0
2	H	126	SER	2.0
1	A	83	ASP	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.