



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 11:18 am BST

PDB ID : 5ZL2
Title : Crystal structure of Bourbon virus envelope glycoprotein at pH8.0
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Deposited on : 2018-03-26
Resolution : 2.70 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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 : 2.11
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) : 2.11

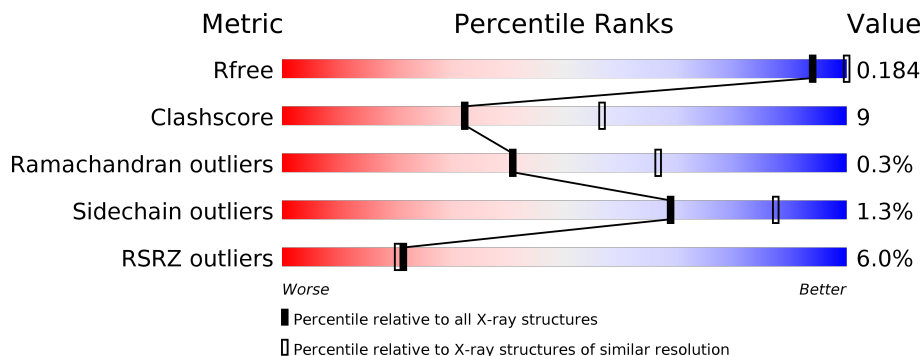
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	467	 % 69% 14% 17%
1	B	467	 6% 63% 19% 17%
1	C	467	 7% 65% 18% 17%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9383 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 glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	388	3089	1960	528	580	21	0	0	0
1	B	388	3089	1960	528	580	21	0	0	0
1	C	388	3089	1960	528	580	21	0	0	0

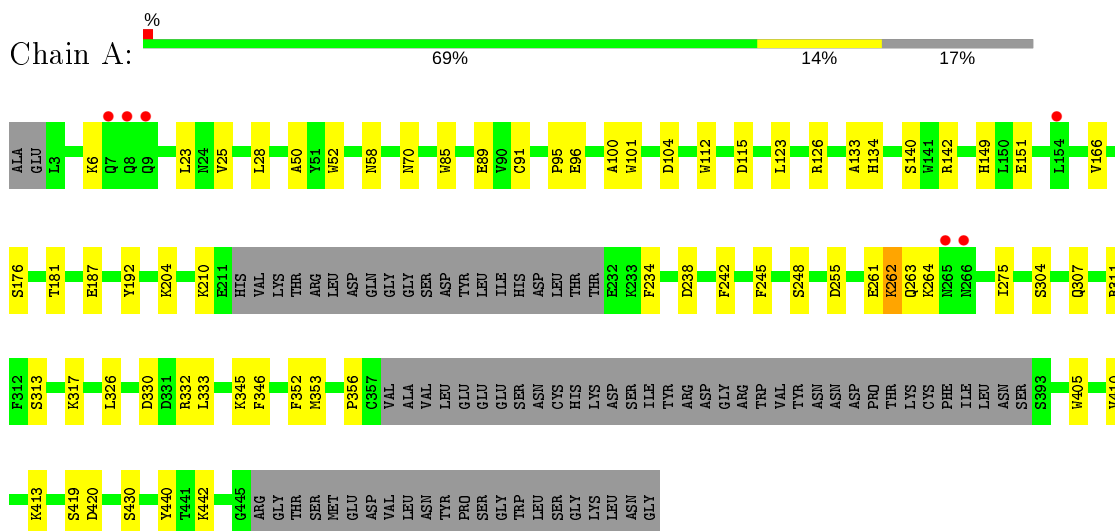
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	31	Total 31	O 31	0	0
2	B	47	Total 47	O 47	0	0
2	C	38	Total 38	O 38	0	0

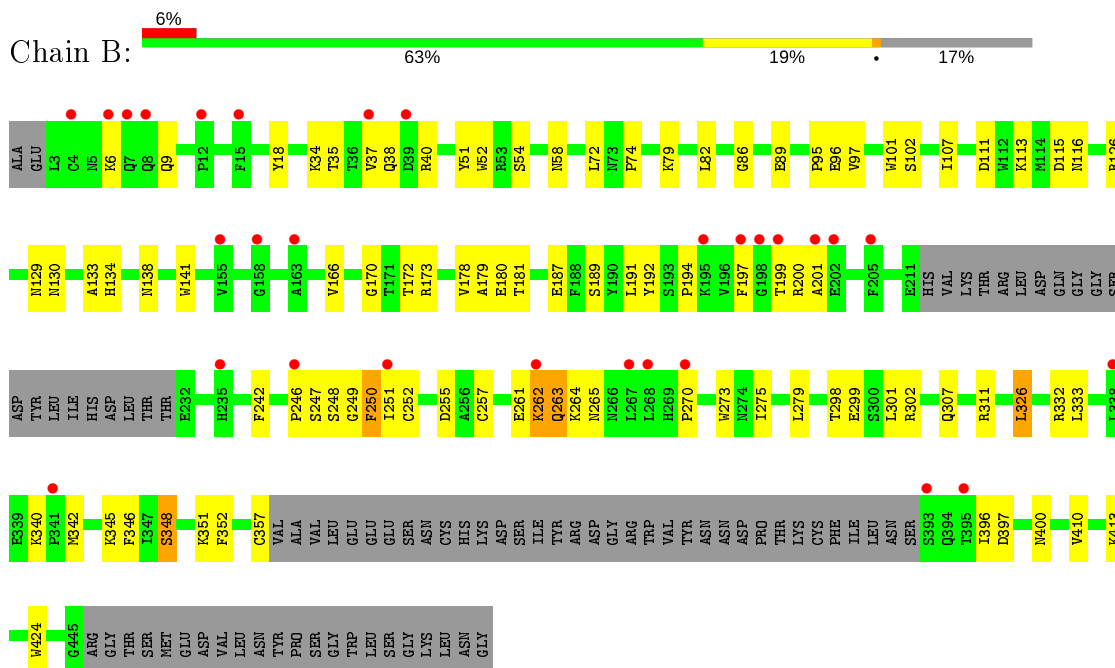
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

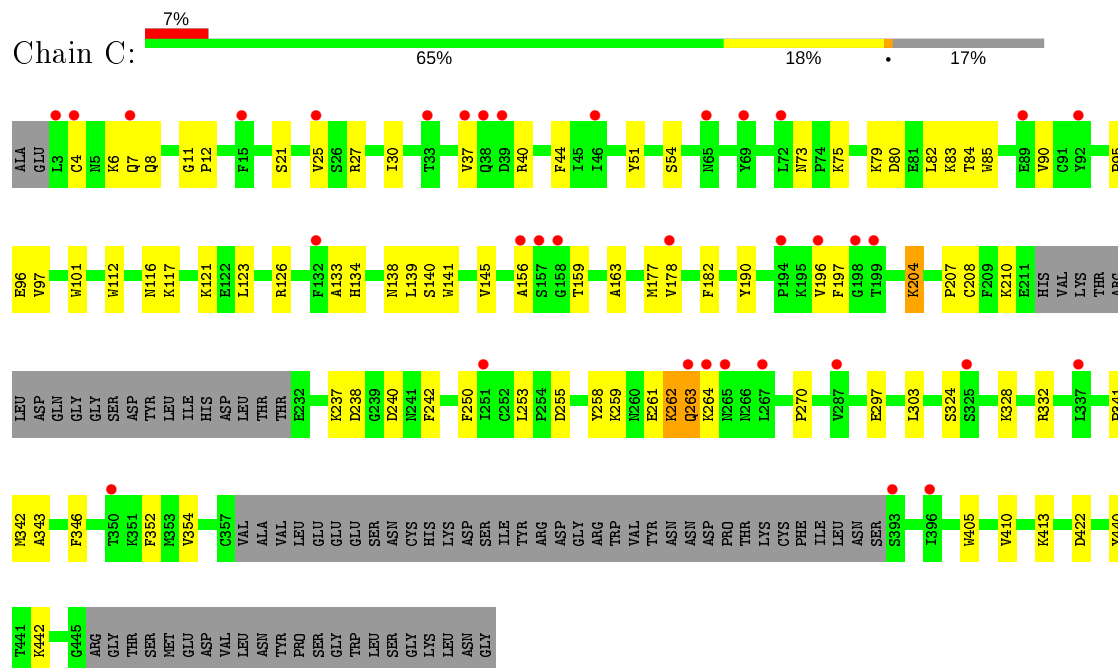
- Molecule 1: Envelope glycoprotein



- Molecule 1: Envelope glycoprotein



- Molecule 1: Envelope glycoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	102.25Å 102.25Å 134.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.79 – 2.70 47.79 – 2.70	Depositor EDS
% Data completeness (in resolution range)	80.6 (47.79-2.70) 80.6 (47.79-2.70)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.201 , 0.240 0.160 , 0.184	Depositor DCC
R_{free} test set	1676 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 16.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.32$, $\langle L^2 \rangle = 0.15$	Xtriage
Estimated twinning fraction	0.399 for -h,-k,l 0.417 for h,-h-k,-l 0.427 for -k,-h,-l	Xtriage
Reported twinning fraction	0.480 for -k,-h,-l	Depositor
Outliers	0 of 34756 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9383	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.18% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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	0.26	0/3165	0.46	0/4284
1	B	0.27	0/3165	0.46	0/4284
1	C	0.29	0/3165	0.45	0/4284
All	All	0.28	0/9495	0.46	0/12852

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	A	0	2
1	B	0	2
1	C	0	2
All	All	0	6

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	101	TRP	Peptide
1	A	262	LYS	Peptide
1	B	101	TRP	Peptide
1	B	262	LYS	Peptide
1	C	101	TRP	Peptide
1	C	262	LYS	Peptide

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	3089	0	2996	36	2
1	B	3089	0	2996	67	0
1	C	3089	0	2996	57	0
2	A	31	0	0	4	1
2	B	47	0	0	12	0
2	C	38	0	0	11	0
All	All	9383	0	8988	160	2

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 9.

All (160) 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:90:VAL:HG11	1:C:117:LYS:HG3	1.68	0.76
1:C:440:TYR:O	1:C:442:LYS:NZ	2.23	0.72
1:A:166:VAL:HG11	1:A:275:ILE:HD11	1.72	0.71
1:B:173:ARG:NH2	2:B:508:HOH:O	2.24	0.71
1:B:37:VAL:HG13	1:B:199:THR:HG23	1.71	0.71
1:B:326:LEU:CD2	1:B:333:LEU:HD22	2.21	0.71
1:B:255:ASP:OD2	1:B:413:LYS:NZ	2.24	0.70
1:B:249:GLY:N	2:B:507:HOH:O	2.23	0.70
1:B:166:VAL:O	2:B:501:HOH:O	2.10	0.69
1:B:130:ASN:OD1	2:B:502:HOH:O	2.11	0.68
1:C:7:GLN:NE2	1:C:8:GLN:OE1	2.27	0.67
1:A:330:ASP:OD1	2:A:501:HOH:O	2.13	0.67
1:C:255:ASP:OD2	1:C:413:LYS:NZ	2.27	0.67
1:A:242:PHE:CZ	1:A:410:VAL:HG13	2.31	0.65
1:B:40:ARG:O	2:B:503:HOH:O	2.14	0.65
1:C:324:SER:O	1:C:328:LYS:NZ	2.30	0.65
1:B:424:TRP:NE1	2:B:501:HOH:O	2.30	0.64
1:C:422:ASP:OD2	2:C:501:HOH:O	2.15	0.63
1:C:346:PHE:HA	1:C:352:PHE:HA	1.79	0.62
1:B:326:LEU:HD22	1:B:333:LEU:HD22	1.81	0.62
1:B:178:VAL:HG23	1:B:179:ALA:H	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:297:GLU:OE1	2:C:502:HOH:O	2.16	0.60
1:A:52:TRP:CH2	1:A:70:ASN:HB3	2.37	0.60
1:C:261:GLU:HA	1:C:264:LYS:HB2	1.84	0.59
1:A:326:LEU:HD23	1:A:333:LEU:HD22	1.84	0.59
1:B:180:GLU:HB3	1:B:191:LEU:HD11	1.85	0.59
1:A:133:ALA:O	1:A:134:HIS:ND1	2.35	0.59
1:A:420:ASP:O	2:A:502:HOH:O	2.16	0.58
1:A:58:ASN:O	1:A:133:ALA:N	2.36	0.58
1:C:90:VAL:CG1	1:C:117:LYS:HG3	2.33	0.58
1:A:23:LEU:HD23	1:A:25:VAL:H	1.69	0.58
1:B:54:SER:O	2:B:504:HOH:O	2.17	0.58
1:C:242:PHE:CZ	1:C:410:VAL:HG13	2.38	0.57
1:A:204:LYS:NZ	1:A:238:ASP:OD1	2.26	0.57
1:C:332:ARG:HH22	1:C:341:PRO:HA	1.69	0.57
1:B:326:LEU:HD23	1:B:333:LEU:HD22	1.87	0.56
1:B:129:ASN:ND2	2:B:509:HOH:O	2.38	0.56
1:C:96:GLU:OE2	1:C:126:ARG:NH2	2.37	0.55
1:A:346:PHE:HA	1:A:352:PHE:HA	1.88	0.55
1:C:25:VAL:HA	1:C:210:LYS:O	2.06	0.55
1:B:181:THR:HG23	1:B:192:TYR:HB3	1.87	0.55
1:B:178:VAL:HG23	1:B:179:ALA:N	2.22	0.55
1:B:6:LYS:O	1:B:9:GLN:NE2	2.40	0.55
1:B:332:ARG:NH2	1:B:357:CYS:SG	2.75	0.54
1:B:299:GLU:HA	1:B:302:ARG:HD3	1.89	0.53
1:B:242:PHE:CZ	1:B:410:VAL:HG13	2.43	0.53
1:A:345:LYS:O	1:A:353:MET:N	2.40	0.53
1:A:85:TRP:O	1:A:89:GLU:N	2.41	0.53
1:B:273:TRP:CZ3	1:B:275:ILE:HA	2.43	0.53
1:C:159:THR:O	2:C:503:HOH:O	2.19	0.53
1:C:405:TRP:NE1	2:C:509:HOH:O	2.33	0.52
1:C:121:LYS:NZ	1:C:141:TRP:O	2.43	0.51
1:A:115:ASP:N	1:A:115:ASP:OD1	2.43	0.51
1:B:250:PHE:CD2	1:B:270:PRO:HB3	2.47	0.50
1:B:250:PHE:CG	1:B:270:PRO:HB3	2.46	0.50
1:A:181:THR:HG23	1:A:192:TYR:HB3	1.93	0.50
1:B:52:TRP:HE3	1:B:72:LEU:HG	1.76	0.50
1:A:440:TYR:CZ	1:A:442:LYS:HE3	2.46	0.49
1:B:58:ASN:O	1:B:133:ALA:N	2.45	0.49
1:B:246:PRO:HG3	1:B:251:ILE:HB	1.94	0.48
1:B:35:THR:HG22	1:B:199:THR:HG21	1.94	0.48
1:C:80:ASP:O	1:C:84:THR:HG22	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:VAL:CG1	1:B:199:THR:HG23	2.42	0.48
1:C:332:ARG:HH12	1:C:342:MET:N	2.12	0.48
1:A:50:ALA:HB2	1:A:187:GLU:HG2	1.96	0.48
1:B:74:PRO:HB3	1:B:187:GLU:HG2	1.96	0.48
1:A:149:HIS:NE2	1:A:151:GLU:OE1	2.35	0.48
1:B:348:SER:HB3	1:B:351:LYS:HG2	1.96	0.48
1:B:346:PHE:HA	1:B:352:PHE:HA	1.94	0.48
1:B:199:THR:HG22	1:B:200:ARG:H	1.78	0.47
1:C:156:ALA:O	2:C:503:HOH:O	2.20	0.47
1:A:261:GLU:HA	1:A:264:LYS:HB2	1.97	0.47
1:A:142:ARG:NH1	1:A:187:GLU:OE2	2.46	0.47
1:C:204:LYS:NZ	1:C:238:ASP:OD1	2.43	0.47
1:C:82:LEU:HD23	2:C:514:HOH:O	2.15	0.47
1:B:97:VAL:HG11	1:B:107:ILE:HG21	1.97	0.47
1:C:253:LEU:HD12	1:C:258:TYR:HE1	1.80	0.47
1:A:25:VAL:HA	1:A:210:LYS:O	2.15	0.46
1:A:96:GLU:HG3	1:A:126:ARG:HH22	1.79	0.46
1:C:79:LYS:HG2	1:C:83:LYS:HE3	1.96	0.46
1:C:27:ARG:HH22	1:C:207:PRO:CB	2.28	0.46
1:A:261:GLU:HG3	1:A:264:LYS:HB2	1.98	0.46
1:C:250:PHE:CG	1:C:270:PRO:HB3	2.51	0.46
1:B:34:LYS:O	1:B:201:ALA:HA	2.16	0.46
1:B:95:PRO:HA	1:B:116:ASN:ND2	2.31	0.46
1:C:95:PRO:HA	1:C:116:ASN:ND2	2.31	0.46
1:A:100:ALA:HB1	1:A:104:ASP:HB2	1.98	0.46
1:B:340:LYS:O	1:B:342:MET:HG2	2.16	0.46
1:B:96:GLU:HG3	1:B:126:ARG:HH22	1.81	0.45
1:A:28:LEU:HD23	2:A:510:HOH:O	2.16	0.45
1:C:37:VAL:HG11	1:C:163:ALA:HB2	1.98	0.45
1:A:234:PHE:HB2	1:A:245:PHE:HB3	1.98	0.45
1:A:307:GLN:O	1:A:311:ARG:HG3	2.17	0.45
1:C:262:LYS:O	1:C:263:GLN:HB2	2.16	0.45
1:C:240:ASP:HB2	1:C:413:LYS:HB2	1.98	0.45
1:C:21:SER:O	2:C:504:HOH:O	2.21	0.45
1:C:346:PHE:CG	1:C:352:PHE:HB3	2.52	0.44
1:B:170:GLY:HA2	1:B:279:LEU:HD11	1.99	0.44
1:B:51:TYR:HB3	1:B:141:TRP:CE3	2.53	0.44
1:C:54:SER:HB3	1:C:138:ASN:HB2	1.99	0.44
1:C:133:ALA:O	1:C:134:HIS:ND1	2.50	0.44
1:A:332:ARG:NE	1:A:332:ARG:HA	2.33	0.44
1:B:194:PRO:HG2	1:B:197:PHE:CD2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:261:GLU:HA	1:B:264:LYS:HB2	1.99	0.44
1:B:397:ASP:OD1	1:B:400:ASN:HA	2.18	0.44
1:C:123:LEU:HB2	1:C:140:SER:HB3	2.00	0.44
1:A:346:PHE:CD1	1:A:352:PHE:HB3	2.52	0.44
1:B:138:ASN:ND2	2:B:504:HOH:O	2.27	0.44
1:C:303:LEU:HD23	1:C:303:LEU:HA	1.82	0.44
1:C:30:ILE:HD12	1:C:208:CYS:SG	2.58	0.44
1:B:18:TYR:OH	1:B:311:ARG:NH2	2.47	0.44
1:B:86:GLY:HA2	2:B:522:HOH:O	2.18	0.44
1:A:123:LEU:HB2	1:A:140:SER:HB3	2.00	0.43
1:B:82:LEU:HD22	1:B:189:SER:HB3	2.00	0.43
1:C:44:PHE:CZ	1:C:190:TYR:HB3	2.54	0.43
1:C:51:TYR:HB3	1:C:141:TRP:CE3	2.53	0.43
1:C:342:MET:SD	1:C:354:VAL:HG21	2.58	0.43
1:B:247:SER:OG	1:B:248:SER:N	2.51	0.43
1:C:51:TYR:CZ	1:C:73:ASN:HB2	2.53	0.43
1:C:97:VAL:HB	1:C:139:LEU:HB3	2.00	0.43
1:B:262:LYS:O	1:B:263:GLN:HB2	2.18	0.43
1:B:79:LYS:HZ3	1:B:191:LEU:HD13	1.84	0.43
1:C:11:GLY:HA3	1:C:12:PRO:HA	1.90	0.43
1:B:89:GLU:HA	2:B:522:HOH:O	2.17	0.43
1:C:134:HIS:N	2:C:511:HOH:O	2.52	0.43
1:C:346:PHE:CD1	1:C:352:PHE:HB3	2.54	0.43
1:A:419:SER:N	2:A:514:HOH:O	2.52	0.42
1:B:248:SER:O	1:B:263:GLN:HG2	2.20	0.42
1:B:38:GLN:O	1:B:200:ARG:NH2	2.52	0.42
1:B:346:PHE:CG	1:B:352:PHE:HB3	2.54	0.42
1:A:176:SER:H	1:A:181:THR:HB	1.84	0.42
1:C:177:MET:SD	1:C:178:VAL:HG13	2.60	0.42
1:B:111:ASP:O	1:B:113:LYS:HG2	2.19	0.42
1:B:252:CYS:O	1:B:273:TRP:HD1	2.02	0.42
1:C:343:ALA:N	2:C:507:HOH:O	2.52	0.42
1:C:182:PHE:HD1	2:C:514:HOH:O	2.03	0.42
1:C:250:PHE:CD2	1:C:270:PRO:HB3	2.54	0.42
1:B:35:THR:HB	1:B:257:CYS:HB2	2.01	0.42
1:B:396:ILE:HD12	1:B:396:ILE:HA	1.76	0.42
1:C:6:LYS:HE3	1:C:6:LYS:HB3	1.84	0.42
1:C:75:LYS:HD3	1:C:75:LYS:HA	1.87	0.42
1:A:255:ASP:OD2	1:A:413:LYS:NZ	2.53	0.41
1:C:95:PRO:HD3	1:C:112:TRP:HB3	2.02	0.41
1:C:177:MET:H	1:C:177:MET:HG3	1.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:SER:O	1:A:317:LYS:HG3	2.21	0.41
1:C:85:TRP:HB3	1:C:145:VAL:HG21	2.02	0.41
1:B:248:SER:N	2:B:507:HOH:O	2.54	0.41
1:C:197:PHE:CD1	1:C:197:PHE:N	2.89	0.41
1:B:298:THR:O	1:B:301:LEU:HB3	2.20	0.41
1:B:102:SER:HB3	1:B:134:HIS:CG	2.56	0.41
1:B:97:VAL:HG21	1:B:107:ILE:HD13	2.01	0.41
1:B:348:SER:HB3	1:B:351:LYS:CG	2.51	0.41
1:A:261:GLU:O	1:A:262:LYS:HB2	2.21	0.40
1:B:115:ASP:N	1:B:115:ASP:OD1	2.54	0.40
1:B:166:VAL:HG12	1:B:172:THR:HG22	2.03	0.40
1:B:178:VAL:CG2	1:B:179:ALA:H	2.33	0.40
1:C:196:VAL:C	1:C:197:PHE:HD1	2.25	0.40
1:C:4:CYS:O	2:C:505:HOH:O	2.22	0.40
1:A:95:PRO:HD3	1:A:112:TRP:HB3	2.03	0.40
1:B:262:LYS:HB2	1:B:262:LYS:HE3	1.87	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:SER:OG	1:A:405:TRP:O[2_555]	2.15	0.05
1:A:356:PRO:O	2:A:512:HOH:O[2_555]	2.19	0.01

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	382/467 (82%)	362 (95%)	18 (5%)	2 (0%)	29	54
1	B	382/467 (82%)	362 (95%)	19 (5%)	1 (0%)	41	66
1	C	382/467 (82%)	361 (94%)	20 (5%)	1 (0%)	41	66

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1146/1401 (82%)	1085 (95%)	57 (5%)	4 (0%)	41 66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	263	GLN
1	C	263	GLN
1	A	6	LYS
1	A	263	GLN

5.3.2 Protein sidechains [i](#)

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	344/414 (83%)	341 (99%)	3 (1%)	78 92
1	B	344/414 (83%)	338 (98%)	6 (2%)	60 84
1	C	344/414 (83%)	340 (99%)	4 (1%)	71 88
All	All	1032/1242 (83%)	1019 (99%)	13 (1%)	69 87

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	CYS
1	A	248	SER
1	A	430	SER
1	B	250	PHE
1	B	265	ASN
1	B	307	GLN
1	B	326	LEU
1	B	345	LYS
1	B	348	SER
1	C	40	ARG
1	C	204	LYS
1	C	237	LYS

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Mol	Chain	Res	Type
1	C	259	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	128	ASN
1	B	241	ASN
1	B	265	ASN
1	C	7	GLN

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 carbohydrates 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, 95th 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 ²)	Q<0.9
1	A	388/467 (83%)	-0.22	6 (1%) 73 76	8, 32, 68, 146	0
1	B	388/467 (83%)	0.23	29 (7%) 14 12	10, 40, 108, 188	0
1	C	388/467 (83%)	0.29	35 (9%) 9 7	12, 48, 117, 194	0
All	All	1164/1401 (83%)	0.10	70 (6%) 21 20	8, 39, 106, 194	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	198	GLY	11.3
1	C	157	SER	10.4
1	C	264	LYS	9.7
1	B	6	LYS	8.9
1	C	265	ASN	8.2
1	C	158	GLY	7.4
1	B	267	LEU	7.2
1	C	178	VAL	6.8
1	C	287	VAL	6.2
1	B	37	VAL	6.2
1	C	198	GLY	6.2
1	C	199	THR	6.0
1	C	7	GLN	6.0
1	B	12	PRO	5.8
1	C	393	SER	5.5
1	B	199	THR	5.5
1	B	39	ASP	4.9
1	C	263	GLN	4.7
1	A	8	GLN	4.6
1	B	158	GLY	4.5
1	A	265	ASN	4.5
1	B	8	GLN	4.4
1	B	338	LEU	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	4	CYS	4.3
1	B	268	LEU	4.1
1	C	39	ASP	4.1
1	B	163	ALA	4.1
1	C	4	CYS	4.1
1	B	197	PHE	3.9
1	C	65	ASN	3.8
1	C	15	PHE	3.7
1	A	9	GLN	3.7
1	B	251	ILE	3.7
1	C	3	LEU	3.6
1	B	235	HIS	3.4
1	C	196	VAL	3.2
1	C	396	ILE	3.2
1	A	7	GLN	3.2
1	B	7	GLN	3.2
1	B	270	PRO	3.0
1	B	15	PHE	2.9
1	B	246	PRO	2.9
1	B	395	THR	2.7
1	A	266	ASN	2.7
1	C	267	LEU	2.7
1	B	195	LYS	2.6
1	B	205	PHE	2.6
1	A	154	LEU	2.5
1	C	337	LEU	2.5
1	C	46	ILE	2.5
1	C	72	LEU	2.5
1	B	393	SER	2.5
1	C	251	ILE	2.4
1	B	341	PRO	2.4
1	C	33	THR	2.4
1	B	155	VAL	2.3
1	B	201	ALA	2.3
1	C	89	GLU	2.3
1	C	156	ALA	2.3
1	B	202	GLU	2.3
1	C	92	TYR	2.2
1	C	350	THR	2.2
1	C	325	SER	2.2
1	C	37	VAL	2.2
1	C	38	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	194	PRO	2.1
1	B	262	LYS	2.1
1	C	25	VAL	2.1
1	C	69	TYR	2.0
1	C	132	PHE	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 carbohydrates 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.