



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:03 AM UTC

PDB ID : 6LGX / pdb_00006lgx
Title : Structure of Rabies virus glycoprotein at basic pH
Authors : Yang, F.L.; Lin, S.; Ye, F.; Yang, J.; Qi, J.X.; Chen, Z.J.; Lin, X.; Wang, J.C.; Yue, D.; Cheng, Y.W.; Chen, Z.M.; Chen, H.; You, Y.; Zhang, Z.L.; Yang, Y.; Yang, M.; Sun, H.L.; Li, Y.H.; Cao, Y.; Yang, S.Y.; Wei, Y.Q.; Gao, G.F.; Lu, G.W.
Deposited on : 2019-12-06
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

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>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

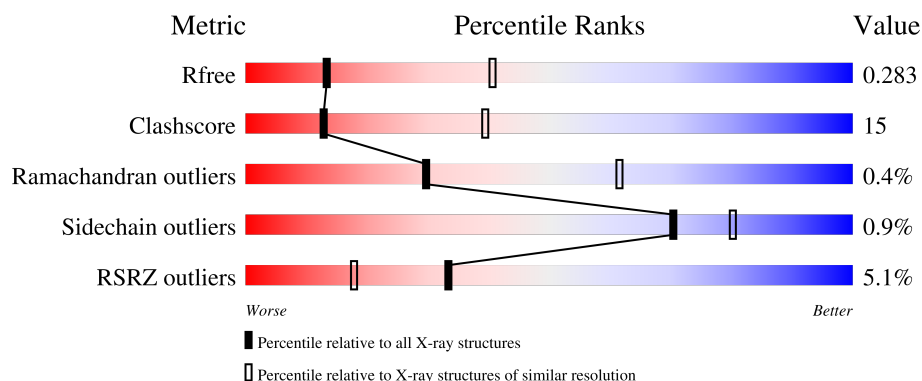
1 Overall quality at a glance

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}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (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 ≥ 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	443	
1	B	443	

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 5695 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 Glycoprotein,Glycoprotein,Glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	377	Total	C	N	O	S	0	0	0
			2968	1881	512	551	24			
1	B	348	Total	C	N	O	S	0	0	0
			2727	1731	473	499	24			

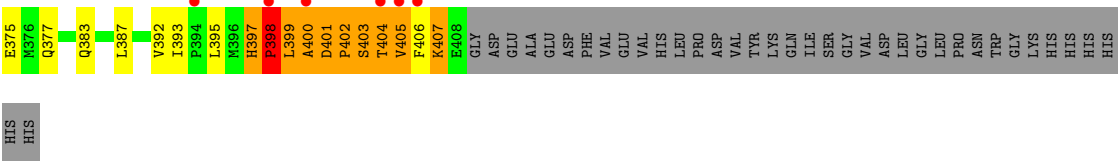
There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ALA	-	expression tag	UNP Q5JZZ2
A	-2	ASP	-	expression tag	UNP Q5JZZ2
A	-1	GLU	-	expression tag	UNP Q5JZZ2
A	0	PHE	-	expression tag	UNP Q5JZZ2
A	75	GLY	-	linker	UNP Q5JZZ2
A	76	GLY	-	linker	UNP Q5JZZ2
A	77	SER	-	linker	UNP Q5JZZ2
A	78	GLY	-	linker	UNP Q5JZZ2
A	79	GLY	-	linker	UNP Q5JZZ2
A	121	GLY	-	linker	UNP Q2Z2I1
A	122	GLY	-	linker	UNP Q2Z2I1
A	123	SER	-	linker	UNP Q2Z2I1
A	124	GLY	-	linker	UNP Q2Z2I1
A	125	GLY	-	linker	UNP Q2Z2I1
A	440	HIS	-	expression tag	UNP D8VEC1
A	441	HIS	-	expression tag	UNP D8VEC1
A	442	HIS	-	expression tag	UNP D8VEC1
A	443	HIS	-	expression tag	UNP D8VEC1
A	444	HIS	-	expression tag	UNP D8VEC1
A	445	HIS	-	expression tag	UNP D8VEC1
B	-3	ALA	-	expression tag	UNP Q5JZZ2
B	-2	ASP	-	expression tag	UNP Q5JZZ2
B	-1	GLU	-	expression tag	UNP Q5JZZ2
B	0	PHE	-	expression tag	UNP Q5JZZ2
B	75	GLY	-	linker	UNP Q5JZZ2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	76	GLY	-	linker	UNP Q5JZZ2
B	77	SER	-	linker	UNP Q5JZZ2
B	78	GLY	-	linker	UNP Q5JZZ2
B	79	GLY	-	linker	UNP Q5JZZ2
B	121	GLY	-	linker	UNP Q2Z2I1
B	122	GLY	-	linker	UNP Q2Z2I1
B	123	SER	-	linker	UNP Q2Z2I1
B	124	GLY	-	linker	UNP Q2Z2I1
B	125	GLY	-	linker	UNP Q2Z2I1
B	440	HIS	-	expression tag	UNP D8VEC1
B	441	HIS	-	expression tag	UNP D8VEC1
B	442	HIS	-	expression tag	UNP D8VEC1
B	443	HIS	-	expression tag	UNP D8VEC1
B	444	HIS	-	expression tag	UNP D8VEC1
B	445	HIS	-	expression tag	UNP D8VEC1



HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	192.44Å 138.45Å 61.44Å 90.00° 100.02° 90.00°	Depositor
Resolution (Å)	47.40 – 3.10 47.40 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.40-3.10) 99.1 (47.40-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.16-3549	Depositor
R, R_{free}	0.242 , 0.285 0.246 , 0.283	Depositor DCC
R_{free} test set	1348 reflections (4.66%)	wwPDB-VP
Wilson B-factor (Å ²)	107.1	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 65.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5695	wwPDB-VP
Average B, all atoms (Å ²)	123.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0/3041	0.48	1/4118 (0.0%)
1	B	0.32	0/2796	0.69	13/3788 (0.3%)
All	All	0.29	0/5837	0.59	14/7906 (0.2%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	158	LYS	N-CA-C	12.21	124.59	111.28
1	B	399	LEU	N-CA-C	11.40	123.71	111.28
1	B	199	ARG	N-CA-C	8.90	120.75	111.14
1	B	200	ALA	N-CA-C	7.56	121.94	111.74
1	A	407	LYS	N-CA-C	7.27	120.24	110.35
1	B	404	THR	N-CA-C	6.05	120.36	112.92
1	B	407	LYS	N-CA-C	5.73	119.47	111.74
1	B	398	PRO	CA-N-CD	-5.67	104.06	112.00
1	B	400	ALA	N-CA-C	5.61	118.16	111.71
1	B	402	PRO	N-CA-C	5.53	123.85	112.47
1	B	146	ASP	N-CA-C	5.50	120.85	112.54
1	B	31	GLU	N-CA-C	5.17	118.72	111.74
1	B	405	VAL	N-CA-C	5.12	117.87	109.78
1	B	157	GLY	N-CA-C	-5.11	104.57	113.76

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	2968	0	2927	78	1
1	B	2727	0	2695	93	1
All	All	5695	0	5622	171	1

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 15.

All (171) 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:B:35:CYS:SG	1:B:200:ALA:O	1.95	1.24
1:B:155:PRO:O	1:B:158:LYS:HB2	1.58	1.01
1:B:148:SER:HA	1:B:160:SER:O	1.59	1.00
1:B:397:HIS:CG	1:B:398:PRO:HD3	1.98	0.99
1:A:333:ARG:NH2	1:A:334:THR:CG2	2.30	0.94
1:B:32:ASP:OD1	1:B:33:GLU:N	2.04	0.91
1:A:333:ARG:CZ	1:A:334:THR:HG23	2.03	0.88
1:A:333:ARG:NH2	1:A:334:THR:HG21	1.89	0.87
1:A:27:ASN:OD1	1:A:276:LEU:CD2	2.23	0.86
1:A:26:ASN:OD1	1:A:28:LEU:HB2	1.74	0.85
1:A:27:ASN:HD22	1:A:308:VAL:HG21	1.46	0.80
1:B:145:TYR:OH	1:B:252:CYS:O	1.99	0.80
1:A:333:ARG:CZ	1:A:334:THR:CG2	2.61	0.79
1:B:406:PHE:O	1:B:407:LYS:HG2	1.84	0.77
1:B:402:PRO:O	1:B:403:SER:HB2	1.86	0.76
1:A:233:LEU:HD11	1:A:246:SER:HB3	1.67	0.76
1:A:27:ASN:HD22	1:A:308:VAL:CG2	2.00	0.74
1:A:231:LEU:O	1:A:244:GLN:NE2	2.20	0.74
1:B:35:CYS:SG	1:B:201:SER:C	2.70	0.74
1:B:403:SER:OG	1:B:404:THR:N	2.20	0.73
1:B:8:PRO:HB3	1:B:327:ALA:HB1	1.71	0.72
1:B:397:HIS:H	1:B:397:HIS:CD2	2.08	0.71
1:B:146:ASP:OD2	1:B:148:SER:HB2	1.91	0.70
1:A:27:ASN:ND2	1:A:308:VAL:HG21	2.06	0.69
1:B:397:HIS:CD2	1:B:398:PRO:HD3	2.28	0.69
1:A:231:LEU:HD12	1:A:244:GLN:HG3	1.74	0.68
1:B:397:HIS:ND1	1:B:398:PRO:HD3	2.08	0.67
1:B:221:GLY:HA3	1:B:236:MET:HE2	1.77	0.67
1:A:27:ASN:ND2	1:A:308:VAL:CG2	2.58	0.67
1:A:17:ILE:HD13	1:A:324:GLU:HB3	1.78	0.66
1:A:105:ASP:O	1:A:107:ARG:NH1	2.29	0.65
1:A:319:ASN:O	1:A:320:LYS:HB2	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:HIS:HD1	1:A:176:THR:HG1	1.42	0.64
1:A:85:LYS:HG3	1:A:86:HIS:H	1.62	0.64
1:B:162:ILE:O	1:B:162:ILE:HG13	1.96	0.64
1:A:410:ASP:N	1:A:410:ASP:OD1	2.32	0.63
1:B:148:SER:O	1:B:149:LEU:HD23	1.97	0.63
1:A:27:ASN:OD1	1:A:276:LEU:HD23	1.98	0.62
1:B:402:PRO:O	1:B:403:SER:CB	2.45	0.62
1:A:333:ARG:HH22	1:A:334:THR:HG21	1.63	0.62
1:B:400:ALA:HB3	1:B:402:PRO:HD3	1.82	0.62
1:B:11:LEU:HD23	1:B:347:VAL:HG11	1.82	0.61
1:B:341:SER:OG	1:B:342:LYS:N	2.33	0.61
1:B:364:ILE:HD11	1:B:372:LEU:HD23	1.81	0.61
1:A:62:THR:OG1	1:A:133:ILE:HG23	2.00	0.61
1:B:142:LEU:HD11	1:B:147:LYS:HA	1.83	0.61
1:B:100:TRP:HE3	1:B:105:ASP:HB3	1.66	0.61
1:B:148:SER:CA	1:B:160:SER:O	2.45	0.61
1:B:338:ILE:HG22	1:B:339:ILE:HG13	1.82	0.61
1:A:201:SER:O	1:A:202:ASN:HB3	2.01	0.60
1:A:2:PHE:CZ	1:A:27:ASN:OD1	2.54	0.59
1:A:405:VAL:HG23	1:A:405:VAL:O	2.03	0.59
1:B:234:ARG:NH1	1:B:238:GLY:O	2.36	0.59
1:B:146:ASP:O	1:B:147:LYS:HB2	2.02	0.59
1:B:31:GLU:HG2	1:B:31:GLU:O	2.02	0.58
1:A:27:ASN:HD21	1:A:308:VAL:HG11	1.67	0.58
1:B:100:TRP:CD1	1:B:107:ARG:HE	2.21	0.58
1:A:281:GLU:HA	1:A:284:LEU:HD12	1.86	0.57
1:B:37:ASN:O	1:B:200:ALA:HA	2.03	0.57
1:A:67:GLU:HB2	1:A:83:LYS:HG3	1.85	0.57
1:B:368:ASP:N	1:B:368:ASP:OD1	2.38	0.56
1:B:397:HIS:CD2	1:B:397:HIS:N	2.73	0.56
1:B:282:GLU:OE2	1:B:282:GLU:N	2.33	0.56
1:B:354:HIS:HB3	1:B:359:PHE:CE1	2.41	0.56
1:B:37:ASN:C	1:B:38:LEU:HD22	2.31	0.56
1:A:282:GLU:OE1	1:A:282:GLU:N	2.31	0.56
1:A:358:VAL:HG12	1:A:364:ILE:HG22	1.87	0.56
1:A:224:ARG:HH21	1:A:248:GLU:HA	1.71	0.55
1:B:407:LYS:O	1:B:407:LYS:HG3	2.05	0.55
1:A:27:ASN:OD1	1:A:276:LEU:HD22	2.03	0.55
1:B:101:LYS:HD3	1:B:134:ILE:HG12	1.88	0.54
1:B:346:LYS:HE2	1:B:349:GLY:HA2	1.89	0.54
1:A:155:PRO:HG2	1:A:158:LYS:HD2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:SER:HA	1:A:206:THR:HA	1.90	0.54
1:B:397:HIS:CE1	1:B:398:PRO:HD3	2.43	0.54
1:A:202:ASN:CG	1:A:203:GLY:H	2.15	0.53
1:B:372:LEU:CD1	1:B:377:GLN:HG3	2.38	0.53
1:B:372:LEU:HD12	1:B:377:GLN:HG3	1.90	0.53
1:B:334:THR:O	1:B:337:GLU:HG2	2.09	0.53
1:A:368:ASP:OD1	1:A:368:ASP:N	2.41	0.53
1:B:332:VAL:HG21	1:B:338:ILE:HD11	1.90	0.53
1:A:363:ILE:HG12	1:A:373:ILE:HG12	1.90	0.52
1:B:288:GLU:O	1:B:292:THR:OG1	2.20	0.51
1:B:405:VAL:O	1:B:407:LYS:NZ	2.41	0.51
1:B:32:ASP:O	1:B:33:GLU:HB2	2.09	0.51
1:B:199:ARG:O	1:B:208:GLY:O	2.28	0.51
1:B:39:SER:O	1:B:198:LYS:HA	2.11	0.51
1:B:62:THR:HG23	1:B:133:ILE:HB	1.93	0.51
1:B:392:VAL:HG12	1:B:393:ILE:H	1.76	0.50
1:A:368:ASP:OD2	1:A:370:HIS:ND1	2.44	0.50
1:B:196:ARG:HB2	1:B:196:ARG:HH11	1.77	0.50
1:A:273:VAL:HG22	1:A:276:LEU:HG	1.93	0.50
1:A:27:ASN:ND2	1:A:308:VAL:CG1	2.75	0.50
1:B:96:ALA:O	1:B:100:TRP:HD1	1.94	0.50
1:B:293:THR:HG23	1:B:295:SER:H	1.76	0.50
1:A:66:THR:HG22	1:A:84:ARG:HB3	1.94	0.49
1:A:202:ASN:CG	1:A:203:GLY:N	2.70	0.49
1:A:2:PHE:HZ	1:A:27:ASN:OD1	1.94	0.49
1:A:224:ARG:NH2	1:A:248:GLU:HG2	2.28	0.49
1:A:8:PRO:O	1:A:347:VAL:HG23	2.13	0.48
1:A:347:VAL:HG12	1:A:352:HIS:HB2	1.94	0.48
1:A:334:THR:O	1:A:337:GLU:HG2	2.14	0.48
1:B:319:ASN:O	1:B:320:LYS:HG3	2.13	0.48
1:A:60:THR:HB	1:A:135:ILE:HB	1.95	0.48
1:A:375:GLU:OE1	1:A:375:GLU:N	2.43	0.48
1:B:65:VAL:HG13	1:B:85:LYS:HB3	1.96	0.48
1:B:204:ASN:OD1	1:B:204:ASN:N	2.46	0.48
1:B:401:ASP:N	1:B:402:PRO:CD	2.77	0.47
1:A:27:ASN:ND2	1:A:308:VAL:HG11	2.27	0.47
1:A:316:THR:HB	1:A:360:PHE:HA	1.96	0.47
1:B:319:ASN:C	1:B:320:LYS:CG	2.87	0.47
1:B:22:LEU:HD21	1:B:322:LEU:HD21	1.95	0.47
1:B:35:CYS:SG	1:B:201:SER:CA	3.02	0.47
1:B:153:VAL:HG22	1:B:171:THR:HG23	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:397:HIS:CG	1:B:398:PRO:CD	2.85	0.47
1:B:46:LEU:HB2	1:B:51:ILE:HG12	1.96	0.47
1:A:101:LYS:HD2	1:A:106:PRO:HG2	1.97	0.47
1:B:58:GLY:HA3	1:B:178:TRP:CZ2	2.49	0.47
1:B:405:VAL:HG12	1:B:406:PHE:N	2.29	0.47
1:B:155:PRO:O	1:B:158:LYS:CB	2.47	0.46
1:A:27:ASN:ND2	1:A:27:ASN:C	2.71	0.46
1:B:405:VAL:HG12	1:B:407:LYS:H	1.81	0.46
1:A:385:MET:O	1:A:389:LYS:N	2.36	0.46
1:B:286:ALA:O	1:B:290:ILE:HG13	2.15	0.46
1:B:166:SER:C	1:B:168:TYR:H	2.24	0.46
1:B:32:ASP:CG	1:B:33:GLU:H	2.10	0.46
1:A:151:SER:H	1:A:157:GLY:HA2	1.81	0.46
1:A:306:LYS:HG2	1:A:308:VAL:O	2.16	0.46
1:B:97:ALA:O	1:B:101:LYS:HD2	2.14	0.46
1:B:60:THR:O	1:B:134:ILE:HA	2.16	0.45
1:A:17:ILE:HB	1:A:21:HIS:HB2	1.99	0.45
1:A:224:ARG:HH21	1:A:249:THR:H	1.65	0.45
1:B:97:ALA:C	1:B:101:LYS:HD2	2.41	0.45
1:B:392:VAL:HG12	1:B:393:ILE:N	2.32	0.45
1:A:62:THR:HG22	1:A:176:THR:OG1	2.17	0.45
1:A:372:LEU:HD12	1:A:372:LEU:O	2.17	0.45
1:A:66:THR:O	1:A:128:THR:HA	2.17	0.44
1:A:2:PHE:HZ	1:A:276:LEU:HD23	1.83	0.44
1:B:208:GLY:HA2	1:B:219:LEU:HG	1.98	0.44
1:A:286:ALA:O	1:A:290:ILE:HG13	2.18	0.44
1:B:196:ARG:HB2	1:B:196:ARG:NH1	2.33	0.44
1:B:41:PHE:CE2	1:B:43:TYR:HB3	2.53	0.44
1:B:404:THR:O	1:B:404:THR:OG1	2.32	0.44
1:A:95:ARG:HB2	1:A:179:MET:HE1	1.99	0.43
1:B:30:VAL:HG11	1:B:218:SER:HB3	1.99	0.43
1:B:287:LEU:HD11	1:B:291:MET:HE2	2.01	0.43
1:B:398:PRO:HD2	1:B:399:LEU:HD23	2.01	0.43
1:A:38:LEU:HG	1:A:200:ALA:HB2	2.01	0.43
1:A:85:LYS:HG3	1:A:86:HIS:N	2.30	0.43
1:A:184:ARG:HG2	1:A:185:PRO:HD3	2.01	0.43
1:A:333:ARG:CZ	1:A:334:THR:HG21	2.39	0.43
1:A:341:SER:OG	1:A:342:LYS:N	2.51	0.43
1:A:6:THR:OG1	1:A:306:LYS:O	2.34	0.43
1:B:383:GLN:O	1:B:387:LEU:HD23	2.18	0.42
1:A:277:VAL:HG13	1:A:280:ARG:HH21	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:306:LYS:NZ	1:B:311:PHE:O	2.38	0.42
1:B:405:VAL:CG1	1:B:406:PHE:N	2.82	0.42
1:B:35:CYS:SG	1:B:202:ASN:N	2.92	0.42
1:A:342:LYS:HG3	1:A:343:GLY:N	2.33	0.42
1:A:226:LYS:HG2	1:A:231:LEU:HD23	2.01	0.42
1:B:173:HIS:ND1	1:B:176:THR:HG23	2.35	0.42
1:B:19:ILE:O	1:B:291:MET:HE1	2.18	0.41
1:B:289:SER:O	1:B:293:THR:HG22	2.19	0.41
1:B:375:GLU:OE1	1:B:375:GLU:N	2.48	0.41
1:B:280:ARG:O	1:B:284:LEU:HG	2.20	0.41
1:A:43:TYR:HE2	1:A:45:GLU:HB3	1.84	0.41
1:A:58:GLY:HA3	1:A:178:TRP:CZ2	2.56	0.41
1:A:397:HIS:HB3	1:A:400:ALA:HB3	2.02	0.41
1:A:153:VAL:HG23	1:A:154:PHE:CD1	2.56	0.40
1:A:211:ASP:N	1:A:211:ASP:OD1	2.51	0.40
1:B:255:ASP:OD1	1:B:255:ASP:N	2.54	0.40

All (1) 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:320:LYS:O	1:B:395:LEU:CD1[1_556]	1.91	0.29

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	369/443 (83%)	345 (94%)	24 (6%)	0	100	100
1	B	340/443 (77%)	313 (92%)	24 (7%)	3 (1%)	14	44
All	All	709/886 (80%)	658 (93%)	48 (7%)	3 (0%)	30	61

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	401	ASP
1	B	403	SER
1	B	398	PRO

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	336/390 (86%)	335 (100%)	1 (0%)	86	87
1	B	309/390 (79%)	304 (98%)	5 (2%)	55	75
All	All	645/780 (83%)	639 (99%)	6 (1%)	70	80

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	B	159	CYS
1	B	160	SER
1	B	321	THR
1	B	397	HIS
1	B	398	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	99	ASN
1	A	384	HIS
1	B	27	ASN
1	B	303	HIS
1	B	397	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 oligosaccharides 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	377/443 (85%)	0.21	15 (3%)	42 23	61, 117, 176, 218	0
1	B	348/443 (78%)	0.41	22 (6%)	26 14	69, 122, 191, 211	0
All	All	725/886 (81%)	0.31	37 (5%)	33 18	61, 119, 183, 218	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	164	VAL	4.9
1	B	144	PRO	4.8
1	A	27	ASN	4.6
1	B	163	THR	4.5
1	B	65	VAL	4.2
1	B	147	LYS	4.0
1	A	236	MET	3.9
1	B	394	PRO	3.5
1	A	54	ILE	3.4
1	A	85	LYS	3.3
1	B	227	LEU	3.2
1	B	400	ALA	3.2
1	B	404	THR	2.9
1	A	133	ILE	2.8
1	A	146	ASP	2.8
1	B	162	ILE	2.8
1	B	406	PHE	2.8
1	B	405	VAL	2.7
1	A	398	PRO	2.7
1	B	225	LEU	2.7
1	A	254	PRO	2.6
1	A	108	TYR	2.6
1	A	102	MET	2.6
1	A	280	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	64	VAL	2.5
1	B	161	GLY	2.5
1	B	398	PRO	2.5
1	A	136	SER	2.5
1	B	298	PHE	2.4
1	B	132	LEU	2.4
1	B	367	PRO	2.4
1	B	108	TYR	2.4
1	A	141	ASP	2.3
1	A	2	PHE	2.2
1	A	250	LYS	2.1
1	B	369	ASP	2.1
1	B	358	VAL	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 oligosaccharides 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.