



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:23 AM UTC

PDB ID : 8ZPQ / pdb_00008zpq
Title : Crystal structure of SARS-Cov-2-BQ1.1-RBD and 70fab
Authors : Wu, Y.; Gao, F.
Deposited on : 2024-05-31
Resolution : 2.75 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

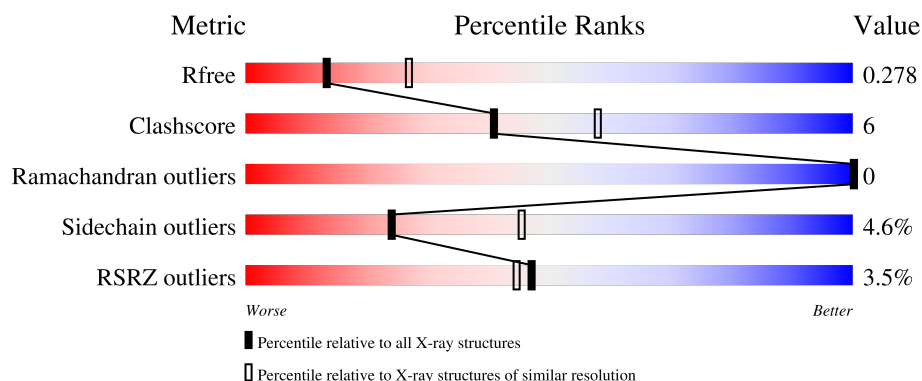
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	228	 4% 72% 21% 5%
1	H	228	 % 75% 21% .
2	B	216	 3% 82% 15% ..
2	L	216	 % 78% 20% .
3	E	223	 4% 71% 17% 12%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	F	223	6% 75% 9% 16%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9519 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 70fab-H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	220	Total	C	N	O	S	0	0	0
			1654	1043	278	326	7			
1	A	216	Total	C	N	O	S	0	0	0
			1630	1031	274	318	7			

- Molecule 2 is a protein called 70fab-L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	213	Total	C	N	O	S	0	0	0
			1591	995	268	324	4			
2	B	212	Total	C	N	O	S	0	0	0
			1583	990	267	322	4			

- Molecule 3 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	197	Total	C	N	O	S	0	0	0
			1566	1009	263	286	8			
3	F	188	Total	C	N	O	S	0	0	0
			1491	956	252	275	8			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	21	ASP	GLY	variant	UNP P0DTC2
E	28	THR	ARG	variant	UNP P0DTC2
E	53	PHE	SER	variant	UNP P0DTC2
E	55	PRO	SER	variant	UNP P0DTC2
E	57	PHE	SER	variant	UNP P0DTC2
E	58	ALA	THR	variant	UNP P0DTC2
E	87	ASN	ASP	variant	UNP P0DTC2
E	90	SER	ARG	variant	UNP P0DTC2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	99	ASN	LYS	variant	UNP P0DTC2
E	122	LYS	ASN	variant	UNP P0DTC2
E	126	THR	LYS	variant	UNP P0DTC2
E	134	ARG	LEU	variant	UNP P0DTC2
E	159	ASN	SER	variant	UNP P0DTC2
E	160	LYS	THR	variant	UNP P0DTC2
E	166	ALA	GLU	variant	UNP P0DTC2
E	168	VAL	PHE	variant	UNP P0DTC2
E	180	ARG	GLN	variant	UNP P0DTC2
E	183	TYR	ASN	variant	UNP P0DTC2
E	187	HIS	TYR	variant	UNP P0DTC2
F	21	ASP	GLY	variant	UNP P0DTC2
F	28	THR	ARG	variant	UNP P0DTC2
F	53	PHE	SER	variant	UNP P0DTC2
F	55	PRO	SER	variant	UNP P0DTC2
F	57	PHE	SER	variant	UNP P0DTC2
F	58	ALA	THR	variant	UNP P0DTC2
F	87	ASN	ASP	variant	UNP P0DTC2
F	90	SER	ARG	variant	UNP P0DTC2
F	99	ASN	LYS	variant	UNP P0DTC2
F	122	LYS	ASN	variant	UNP P0DTC2
F	126	THR	LYS	variant	UNP P0DTC2
F	134	ARG	LEU	variant	UNP P0DTC2
F	159	ASN	SER	variant	UNP P0DTC2
F	160	LYS	THR	variant	UNP P0DTC2
F	166	ALA	GLU	variant	UNP P0DTC2
F	168	VAL	PHE	variant	UNP P0DTC2
F	180	ARG	GLN	variant	UNP P0DTC2
F	183	TYR	ASN	variant	UNP P0DTC2
F	187	HIS	TYR	variant	UNP P0DTC2

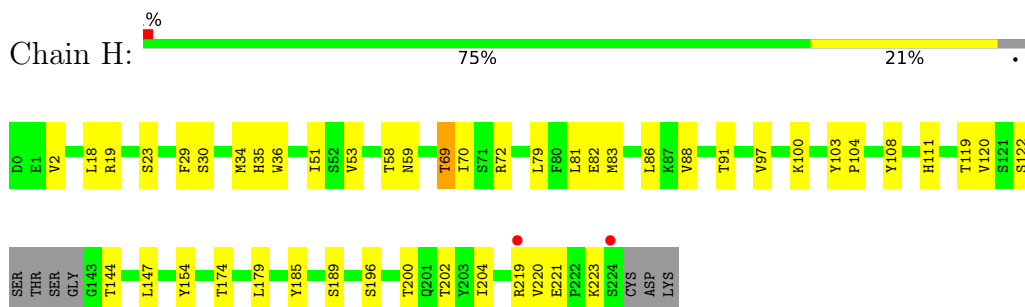
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	H	1	Total O 1 1	0	0
4	L	2	Total O 2 2	0	0
4	A	1	Total O 1 1	0	0

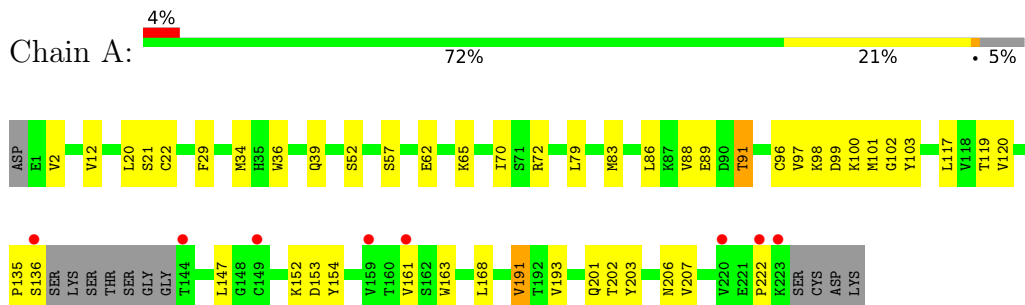
3 Residue-property plots [i](#)

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.

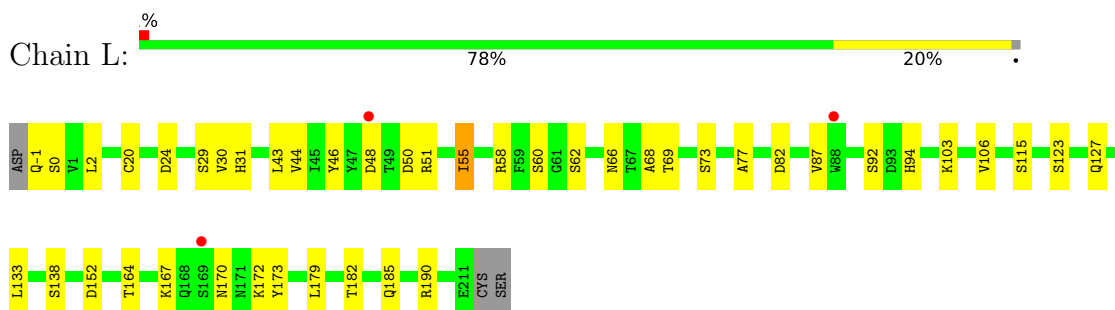
• Molecule 1: 70fab-H



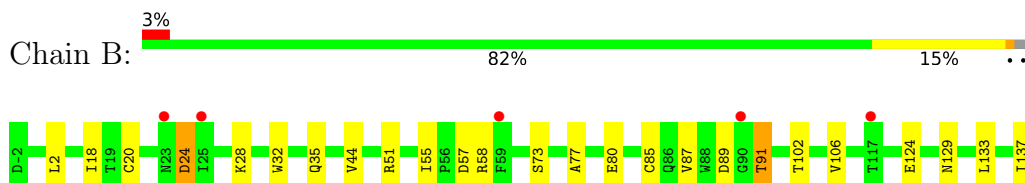
• Molecule 1: 70fab-H



• Molecule 2: 70fab-L

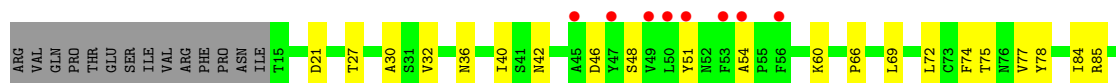


• Molecule 2: 70fab-L

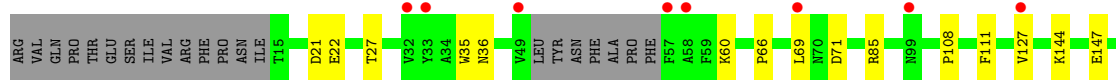
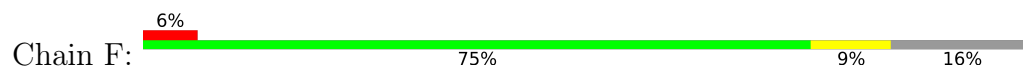




• Molecule 3: Spike protein S1



• Molecule 3: Spike protein S1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.89Å 93.25Å 257.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.35 – 2.75 47.35 – 2.75	Depositor EDS
% Data completeness (in resolution range)	97.4 (47.35-2.75) 97.4 (47.35-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487, PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.234 , 0.281 0.232 , 0.278	Depositor DCC
R_{free} test set	2183 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	66.2	Xtriage
Anisotropy	0.587	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9519	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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

5.1 Standard geometry

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	0.11	0/1669	0.31	0/2270
1	H	0.09	0/1693	0.27	0/2302
2	B	0.12	0/1624	0.29	0/2222
2	L	0.10	0/1632	0.30	0/2233
3	E	0.09	0/1613	0.28	0/2197
3	F	0.13	0/1532	0.33	0/2082
All	All	0.11	0/9763	0.30	0/13306

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

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	1630	0	1603	27	0
1	H	1654	0	1620	22	0
2	B	1583	0	1524	21	0
2	L	1591	0	1533	22	0
3	E	1566	0	1491	24	0
3	F	1491	0	1420	14	0
4	A	1	0	0	0	0
4	H	1	0	0	0	0
4	L	2	0	0	0	0
All	All	9519	0	9191	121	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 (121) 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)
2:B:124:GLU:N	2:B:124:GLU:OE1	2.16	0.78
3:E:143:LEU:HG	3:E:147:GLU:HB3	1.69	0.75
3:F:71:ASP:HA	3:F:211:LYS:HE2	1.69	0.73
2:L:30:VAL:N	2:L:48:ASP:OD1	2.20	0.71
3:F:22:GLU:OE1	3:F:22:GLU:N	2.24	0.68
1:A:91:THR:HG22	1:A:120:VAL:H	1.60	0.67
1:H:91:THR:HG22	1:H:120:VAL:H	1.60	0.66
1:H:108:TYR:HB3	2:L:31:HIS:CD2	2.34	0.63
1:H:29:PHE:O	1:H:72:ARG:NH2	2.32	0.62
3:F:66:PRO:HA	3:F:69:LEU:HG	1.81	0.62
1:H:2:VAL:HG21	1:H:111:HIS:CE1	2.35	0.62
2:L:123:SER:O	2:L:127:GLN:HG2	2.00	0.61
1:A:201:GLN:OE1	1:A:202:THR:N	2.32	0.61
1:A:128:PRO:HB3	1:A:154:TYR:HB3	1.83	0.61
2:L:133:LEU:HB2	2:L:179:LEU:HB3	1.84	0.60
1:A:161:VAL:HG13	1:A:207:VAL:HG12	1.83	0.60
3:E:21:ASP:N	3:E:21:ASP:OD1	2.34	0.60
3:F:108:PRO:HD2	3:F:111:PHE:HB2	1.85	0.59
1:H:174:THR:HG23	1:H:189:SER:HB2	1.85	0.58
3:F:180:ARG:HD3	3:F:183:TYR:HE1	1.68	0.58
1:A:152:LYS:HG2	1:A:153:ASP:OD1	2.03	0.57
3:E:54:ALA:HB3	3:E:118:TRP:HB2	1.86	0.57
3:E:60:LYS:HA	3:F:60:LYS:HA	1.87	0.57
1:A:12:VAL:HG11	1:A:86:LEU:HD13	1.86	0.56
3:E:75:THR:O	3:E:205:THR:OG1	2.24	0.56
3:E:42:ASN:H	3:E:205:THR:HB	1.71	0.55
2:B:28:LYS:NZ	2:B:89:ASP:OD1	2.33	0.55
2:B:137:ILE:HG12	2:B:196:VAL:HG21	1.87	0.55
1:H:18:LEU:HD12	1:H:19:ARG:H	1.70	0.55
2:B:147:VAL:HG22	2:B:196:VAL:HG22	1.89	0.55
1:H:34:MET:HB3	1:H:79:LEU:HD22	1.87	0.55
2:B:18:ILE:HD13	2:B:102:THR:HB	1.88	0.55
2:L:51:ARG:NH1	2:L:55:ILE:O	2.40	0.54
3:E:113:GLY:HA2	3:E:197:PHE:CD2	2.42	0.54
2:L:24:ASP:OD2	3:F:27:THR:HG23	2.08	0.54
2:L:66:ASN:OD1	2:L:66:ASN:N	2.41	0.53
2:L:77:ALA:HA	2:L:106:VAL:HG21	1.90	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:35:HIS:HB2	1:H:97:VAL:HG22	1.91	0.52
3:E:91:GLN:HG2	3:E:101:ALA:HB2	1.91	0.52
1:A:135:PRO:HG3	1:A:147:LEU:HD23	1.91	0.52
1:A:168:LEU:HD21	1:A:191:VAL:HG21	1.92	0.52
1:H:69:THR:HG23	1:H:82:GLU:HB3	1.91	0.51
1:A:88:VAL:O	1:A:91:THR:HG23	2.09	0.51
3:E:69:LEU:HD23	3:E:72:LEU:HD12	1.91	0.51
3:E:78:TYR:HB2	3:E:196:SER:HB2	1.93	0.51
2:L:1:GLN:HE22	2:L:94:HIS:HB2	1.75	0.51
3:E:32:VAL:HG22	3:E:104:ASN:HB3	1.93	0.51
3:E:147:GLU:HG3	1:A:103:TYR:CD2	2.45	0.51
2:B:2:LEU:HD11	2:B:87:VAL:HG13	1.92	0.50
1:A:163:TRP:HB3	1:A:168:LEU:HD23	1.93	0.50
3:E:46:ASP:OD1	3:E:48:SER:N	2.39	0.50
2:L:43:LEU:HD21	2:L:46:TYR:HD2	1.76	0.49
2:B:77:ALA:HA	2:B:106:VAL:HG11	1.94	0.49
1:H:128:PRO:HB3	1:H:154:TYR:HB3	1.95	0.49
1:H:51:ILE:HG13	1:H:58:THR:HG22	1.94	0.48
3:E:154:ILE:HD12	3:E:166:ALA:HB2	1.96	0.48
3:E:27:THR:HG22	2:B:24:ASP:OD2	2.13	0.48
3:E:99:ASN:OD1	3:E:99:ASN:N	2.44	0.48
3:F:21:ASP:OD1	3:F:21:ASP:N	2.42	0.47
1:H:219:ARG:NH1	1:H:221:GLU:HB3	2.30	0.47
3:E:40:ILE:HB	3:E:77:VAL:HB	1.97	0.47
1:A:39:GLN:NE2	2:B:35:GLN:OE1	2.40	0.47
3:E:46:ASP:OD1	3:E:46:ASP:C	2.57	0.46
2:L:46:TYR:CE1	2:L:50:ASP:HB3	2.50	0.46
2:B:133:LEU:HB2	2:B:179:LEU:HB3	1.95	0.46
1:A:98:LYS:HE3	1:A:99:ASP:O	2.16	0.46
2:L:82:ASP:OD2	2:L:103:LYS:HE2	2.15	0.46
1:A:193:VAL:HG11	1:A:203:TYR:CE2	2.50	0.46
1:H:83:MET:HE2	1:H:86:LEU:HD21	1.97	0.45
1:A:62:GLU:OE1	1:A:65:LYS:NZ	2.43	0.45
2:L:167:LYS:HG2	2:L:173:TYR:CE2	2.51	0.45
1:A:91:THR:CG2	1:A:120:VAL:H	2.28	0.45
1:H:36:TRP:HD1	1:H:70:ILE:HD12	1.81	0.45
1:A:22:CYS:HB3	1:A:79:LEU:HB3	1.99	0.45
2:B:58:ARG:HB3	2:B:73:SER:O	2.17	0.45
1:A:99:ASP:OD1	1:A:100:LYS:N	2.48	0.45
1:H:100:LYS:NZ	1:H:104:PRO:O	2.50	0.45
3:F:35:TRP:CD1	3:F:35:TRP:H	2.35	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:30:ALA:HA	2:B:91:THR:HG22	1.99	0.44
1:A:36:TRP:HD1	1:A:70:ILE:HD12	1.82	0.44
2:L:133:LEU:HD12	2:L:179:LEU:HD23	2.00	0.43
3:E:66:PRO:HB2	3:F:66:PRO:HB2	1.99	0.43
3:E:210:LYS:HD2	3:E:210:LYS:HA	1.86	0.43
2:L:2:LEU:HD11	2:L:87:VAL:HG13	2.00	0.43
1:H:204:ILE:HG12	1:H:219:ARG:HA	2.01	0.43
3:F:69:LEU:HD23	3:F:69:LEU:HA	1.65	0.43
2:L:20:CYS:N	2:L:68:ALA:O	2.52	0.43
2:B:44:VAL:HA	2:B:55:ILE:HG13	1.99	0.43
2:B:129:ASN:O	2:B:129:ASN:ND2	2.52	0.43
2:L:44:VAL:HA	2:L:55:ILE:HG12	2.01	0.43
1:A:135:PRO:HD2	1:A:222:PRO:HA	2.01	0.43
1:H:219:ARG:NH1	1:H:221:GLU:OE1	2.52	0.42
1:A:52:SER:HB3	1:A:57:SER:HB2	2.01	0.42
1:H:103:TYR:CD1	3:F:147:GLU:HG3	2.54	0.42
1:A:89:GLU:H	1:A:89:GLU:CD	2.27	0.42
2:B:181:LEU:HD22	2:B:185:GLN:HB3	2.01	0.42
2:L:182:THR:OG1	2:L:185:GLN:HG3	2.19	0.42
3:F:180:ARG:HB2	3:F:183:TYR:CE1	2.55	0.42
1:H:86:LEU:HB3	1:H:120:VAL:HG21	2.02	0.42
3:E:85:ARG:HG3	3:E:177:TYR:CE1	2.54	0.42
2:B:80:GLU:OE2	2:B:106:VAL:N	2.47	0.41
2:L:152:ASP:OD1	2:L:190:ARG:HB3	2.19	0.41
2:L:170:ASN:ND2	2:L:172:LYS:HD2	2.35	0.41
1:H:147:LEU:HB2	1:H:220:VAL:HG11	2.03	0.41
1:A:29:PHE:O	1:A:72:ARG:NH2	2.54	0.41
2:B:2:LEU:HD23	2:B:2:LEU:HA	1.85	0.41
1:A:101:MET:HE3	1:A:102:GLY:H	1.86	0.41
2:B:195:GLN:NE2	2:B:204:GLU:OE2	2.54	0.41
2:L:58:ARG:HB3	2:L:73:SER:O	2.21	0.41
3:E:69:LEU:HD23	3:E:69:LEU:HA	1.87	0.41
1:A:20:LEU:HG	1:A:83:MET:HE3	2.03	0.41
1:H:179:LEU:HD13	1:H:185:TYR:CZ	2.56	0.41
2:L:92:SER:C	2:L:94:HIS:H	2.29	0.41
2:B:51:ARG:CZ	2:B:57:ASP:HA	2.51	0.41
1:A:98:LYS:HE2	1:A:98:LYS:HB3	1.98	0.40
3:E:74:PHE:CD2	3:E:197:PHE:HB3	2.56	0.40
2:B:32:TRP:CZ3	2:B:85:CYS:HB3	2.56	0.40
1:H:36:TRP:CG	1:H:81:LEU:HD22	2.56	0.40
3:F:160:LYS:HD3	3:F:161:PRO:HD2	2.02	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:MET:HE3	1:A:34:MET:HB3	1.79	0.40
2:B:20:CYS:HB2	2:B:32:TRP:CH2	2.57	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	212/228 (93%)	211 (100%)	1 (0%)	0	100	100
1	H	216/228 (95%)	213 (99%)	3 (1%)	0	100	100
2	B	210/216 (97%)	203 (97%)	7 (3%)	0	100	100
2	L	211/216 (98%)	204 (97%)	7 (3%)	0	100	100
3	E	195/223 (87%)	192 (98%)	3 (2%)	0	100	100
3	F	182/223 (82%)	180 (99%)	2 (1%)	0	100	100
All	All	1226/1334 (92%)	1203 (98%)	23 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	185/195 (95%)	174 (94%)	11 (6%)	18	33

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	188/195 (96%)	174 (93%)	14 (7%)	13	24
2	B	177/181 (98%)	173 (98%)	4 (2%)	44	66
2	L	178/181 (98%)	169 (95%)	9 (5%)	21	40
3	E	169/195 (87%)	163 (96%)	6 (4%)	31	55
3	F	162/195 (83%)	157 (97%)	5 (3%)	35	59
All	All	1059/1142 (93%)	1010 (95%)	49 (5%)	24	45

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	23	SER
1	H	30	SER
1	H	53	VAL
1	H	59	ASN
1	H	69	THR
1	H	88	VAL
1	H	119	THR
1	H	122	SER
1	H	137	SER
1	H	144	THR
1	H	196	SER
1	H	200	THR
1	H	202	THR
1	H	223	LYS
2	L	0	SER
2	L	29	SER
2	L	55	ILE
2	L	60	SER
2	L	62	SER
2	L	69	THR
2	L	115	SER
2	L	138	SER
2	L	164	THR
3	E	36	ASN
3	E	51	TYR
3	E	84	ILE
3	E	143	LEU
3	E	151	SER
3	E	182	THR
3	F	36	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	F	85	ARG
3	F	127	VAL
3	F	144	LYS
3	F	160	LYS
1	A	2	VAL
1	A	21	SER
1	A	91	THR
1	A	96	CYS
1	A	97	VAL
1	A	117	LEU
1	A	119	THR
1	A	130	VAL
1	A	136	SER
1	A	191	VAL
1	A	206	ASN
2	B	24	ASP
2	B	91	THR
2	B	154	SER
2	B	201	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	39	GLN
1	H	59	ASN
1	H	201	GLN
2	L	35	GLN
3	E	36	ASN
3	E	175	GLN
3	F	42	ASN
3	F	70	ASN
1	A	180	GLN
2	B	23	ASN
2	B	127	GLN

5.3.3 RNA ⓘ

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(Å²)	Q<0.9	
1	A	216/228 (94%)	0.45	9 (4%)	40	39	45, 62, 109, 122	0
1	H	220/228 (96%)	0.25	3 (1%)	73	73	45, 57, 81, 101	0
2	B	212/216 (98%)	0.51	6 (2%)	55	53	51, 77, 96, 106	0
2	L	213/216 (98%)	0.25	3 (1%)	73	73	45, 60, 77, 106	0
3	E	197/223 (88%)	0.64	10 (5%)	33	33	49, 71, 102, 116	0
3	F	188/223 (84%)	0.84	13 (6%)	23	22	55, 80, 116, 125	1 (0%)
All	All	1246/1334 (93%)	0.48	44 (3%)	47	44	45, 66, 103, 125	1 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	161	PRO	5.3
1	A	222	PRO	3.8
3	E	56	PHE	3.5
1	A	223	LYS	3.1
3	F	166	ALA	3.1
1	A	136	SER	2.9
3	E	50	LEU	2.9
3	F	170	CYS	2.9
3	E	51	TYR	2.8
3	F	57	PHE	2.8
3	F	99	ASN	2.7
3	F	32	VAL	2.7
1	H	137	SER	2.6
3	E	54	ALA	2.6
2	B	117	THR	2.6
1	A	220	VAL	2.6
2	B	90	GLY	2.5
3	E	47	TYR	2.5
3	F	49	VAL	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
3	E	53	PHE	2.4
2	L	88	TRP	2.4
3	E	49	VAL	2.4
2	L	48	ASP	2.4
1	H	224	SER	2.4
3	E	127	VAL	2.3
3	F	127	VAL	2.3
3	F	33	TYR	2.3
3	F	152	THR	2.3
2	B	59	PHE	2.3
3	E	45	ALA	2.2
2	B	23	ASN	2.2
1	A	159	VAL	2.2
1	A	144	THR	2.2
3	F	69	LEU	2.2
3	E	199	LEU	2.2
1	A	149	CYS	2.1
2	B	25	ILE	2.1
1	H	219	ARG	2.1
2	B	209	PRO	2.1
3	F	58	ALA	2.1
1	A	161	VAL	2.0
3	F	179	PHE	2.0
1	A	134	ALA	2.0
2	L	169	SER	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.