



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:29 AM UTC

PDB ID : 7LMC / pdb_00007lmc
Title : Structure of SARS CoV-2 main protease shows simultaneous processing of its N- and C-terminii
Authors : Gajiwala, K.S.; Ferre, R.A.; Liu, W.; Ryan, K.
Deposited on : 2021-02-05
Resolution : 2.98 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

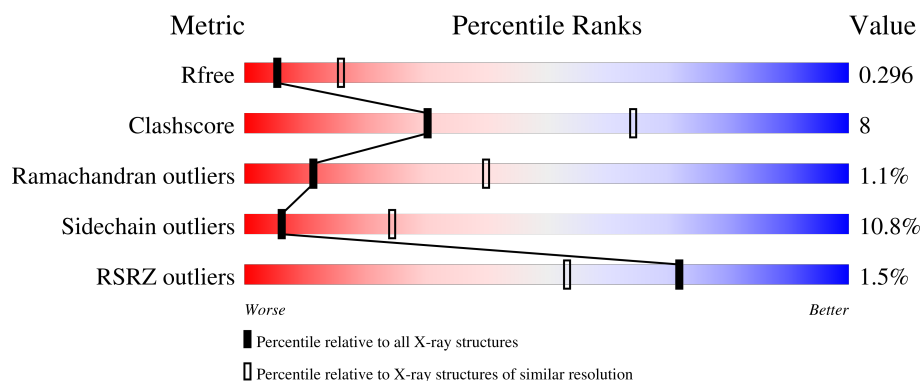
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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	E	6	
1	F	6	
2	A	306	
2	B	306	
2	C	306	

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Mol	Chain	Length	Quality of chain
2	D	306	2% 70% 25%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9539 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 Non-structural protein 4 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	E	6	Total	C	N	O	0	0	0
			42	26	7	9			
1	F	6	Total	C	N	O	0	0	0
			42	26	7	9			

- Molecule 2 is a protein called 3C-like proteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	306	Total	C	N	O	S	0	0	0
			2366	1499	402	444	21			
2	B	306	Total	C	N	O	S	0	0	0
			2366	1499	402	444	21			
2	C	306	Total	C	N	O	S	0	0	0
			2366	1499	402	444	21			
2	D	305	Total	C	N	O	S	0	0	0
			2357	1494	400	442	21			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	145	ALA	CYS	engineered mutation	UNP P0DTD1
B	145	ALA	CYS	engineered mutation	UNP P0DTD1
C	145	ALA	CYS	engineered mutation	UNP P0DTD1
D	145	ALA	CYS	engineered mutation	UNP P0DTD1

3 Residue-property plots

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: Non-structural protein 4 peptide

Chain E: 



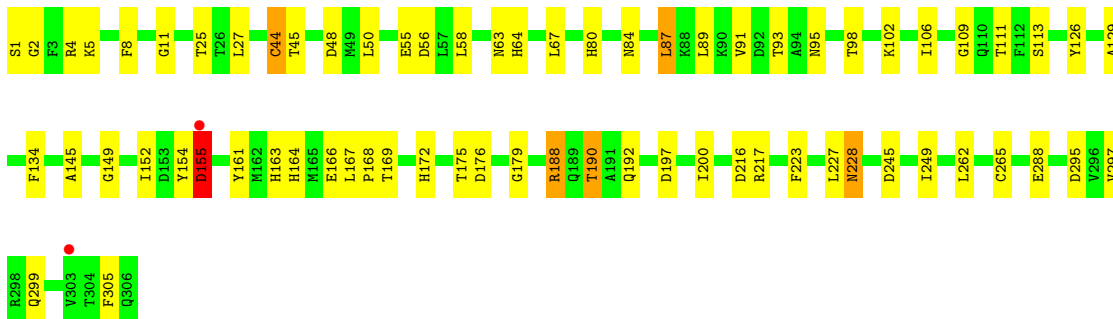
- Molecule 1: Non-structural protein 4 peptide

Chain F: 



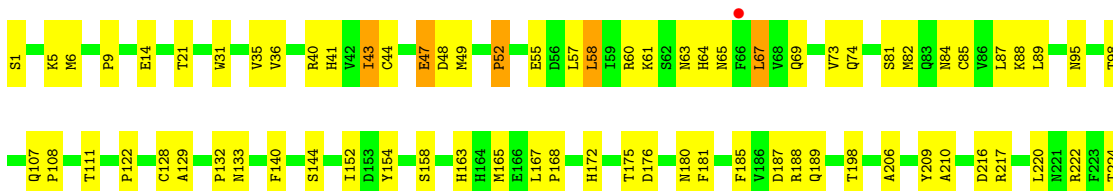
- Molecule 2: 3C-like proteinase

Chain A: 



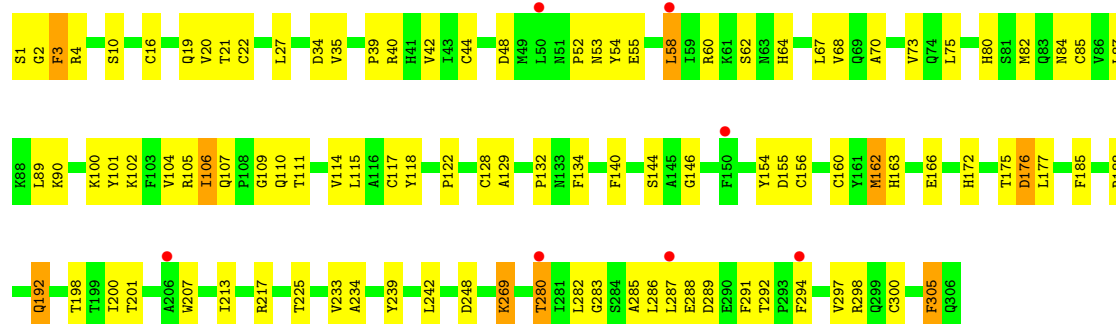
- Molecule 2: 3C-like proteinase

Chain B: 

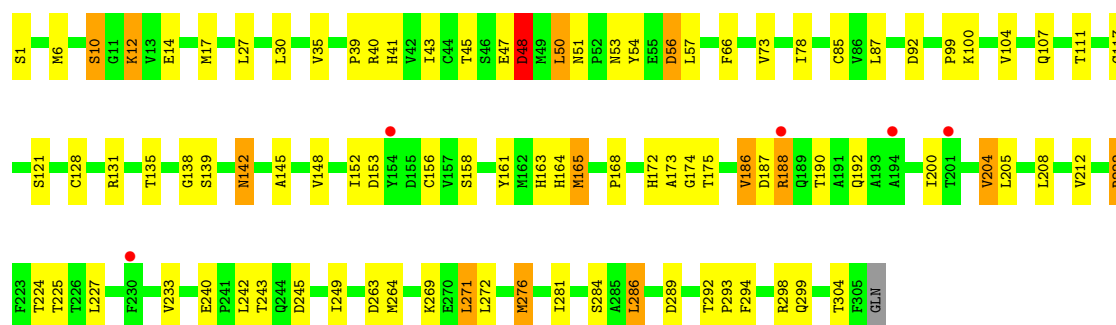




• Molecule 2: 3C-like proteinase



• Molecule 2: 3C-like proteinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.77Å 66.30Å 80.26Å 80.52° 89.46° 70.55°	Depositor
Resolution (Å)	79.07 – 2.98 79.07 – 2.98	Depositor EDS
% Data completeness (in resolution range)	88.7 (79.07-2.98) 88.7 (79.07-2.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.96Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.208 , 0.288 0.217 , 0.296	Depositor DCC
R_{free} test set	1095 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	63.2	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9539	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% 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	E	0.81	0/41	0.92	0/55
1	F	0.62	0/41	1.17	0/55
2	A	0.79	0/2419	1.19	10/3288 (0.3%)
2	B	0.72	0/2419	1.12	5/3288 (0.2%)
2	C	0.77	1/2419 (0.0%)	1.29	5/3288 (0.2%)
2	D	0.78	1/2410 (0.0%)	1.16	4/3276 (0.1%)
All	All	0.77	2/9749 (0.0%)	1.19	24/13250 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	276	MET	SD-CE	-6.14	1.64	1.79
2	C	280	THR	CA-CB	5.03	1.60	1.53

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	134	PHE	CA-CB-CG	-9.53	104.27	113.80
2	C	305	PHE	CA-CB-CG	-8.90	104.90	113.80
2	C	3	PHE	CA-CB-CG	-7.94	105.86	113.80
2	C	176	ASP	CA-CB-CG	7.16	119.76	112.60
2	D	48	ASP	CA-C-N	7.12	130.14	120.38
2	D	48	ASP	C-N-CA	7.12	130.14	120.38
2	A	223	PHE	CA-CB-CG	6.88	120.68	113.80
2	B	176	ASP	CA-CB-CG	6.72	119.32	112.60
2	A	176	ASP	CA-CB-CG	6.11	118.71	112.60
2	B	187	ASP	CA-CB-CG	5.97	118.57	112.60
2	A	228	ASN	CA-CB-CG	-5.92	106.68	112.60
2	C	155	ASP	CA-CB-CG	5.91	118.50	112.60
2	A	63	ASN	CA-C-N	5.73	128.54	120.28
2	A	63	ASN	C-N-CA	5.73	128.54	120.28
2	A	155	ASP	CA-CB-CG	5.72	118.32	112.60
2	A	305	PHE	CA-CB-CG	-5.58	108.22	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	92	ASP	CA-CB-CG	5.45	118.05	112.60
2	A	56	ASP	CA-CB-CG	5.39	117.99	112.60
2	D	187	ASP	CA-CB-CG	5.30	117.90	112.60
2	A	295	ASP	CA-CB-CG	5.13	117.73	112.60
2	C	34	ASP	CA-CB-CG	5.13	117.73	112.60
2	B	133	ASN	CA-C-N	5.08	130.06	122.74
2	B	133	ASN	C-N-CA	5.08	130.06	122.74
2	B	5	LYS	N-CA-C	-5.03	100.31	108.41

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	E	42	0	47	3	0
1	F	42	0	47	7	0
2	A	2366	0	2314	25	0
2	B	2366	0	2314	48	0
2	C	2366	0	2314	43	0
2	D	2357	0	2306	52	0
All	All	9539	0	9342	159	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 8.

All (159) 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:D:51:ASN:HB2	2:D:188:ARG:HE	1.43	0.82
2:B:84:ASN:ND2	2:B:180:ASN:ND2	2.28	0.81
2:B:58:LEU:HD13	2:B:82:MET:HB2	1.63	0.81
2:C:40:ARG:HD3	2:C:85:CYS:HA	1.70	0.73
2:C:285:ALA:H	2:D:286:LEU:HD11	1.53	0.73
2:D:51:ASN:CB	2:D:188:ARG:HE	2.02	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:166:GLU:OE2	2:B:1:SER:HB3	1.88	0.72
2:C:146:GLY:HA2	2:C:162:MET:HE2	1.70	0.72
2:C:213:ILE:HG21	2:C:300:CYS:HB3	1.71	0.71
2:A:188:ARG:HB3	2:A:190:THR:HG22	1.75	0.69
1:E:-1:LEU:HD12	2:B:41:HIS:CG	2.28	0.69
2:D:100:LYS:O	2:D:156:CYS:HA	1.93	0.68
2:A:58:LEU:HD11	2:A:80:HIS:HD2	1.58	0.67
2:A:161:TYR:HH	2:A:163:HIS:HD1	1.43	0.67
2:B:21:THR:HB	2:B:67:LEU:HB3	1.76	0.67
1:F:-1:LEU:HD12	2:D:41:HIS:CG	2.30	0.65
2:C:109:GLY:HA2	2:C:200:ILE:HD13	1.76	0.65
2:B:84:ASN:ND2	2:B:180:ASN:HD21	1.94	0.65
2:A:87:LEU:HD13	2:A:89:LEU:HD21	1.80	0.63
2:C:163:HIS:HE1	2:C:172:HIS:HB3	1.63	0.62
2:D:45:THR:H	2:D:48:ASP:HB2	1.65	0.61
2:C:58:LEU:HD13	2:C:82:MET:HE2	1.83	0.60
2:B:243:THR:HG23	2:B:246:HIS:CE1	2.37	0.60
2:B:111:THR:HG22	2:B:129:ALA:HB2	1.83	0.59
2:C:233:VAL:HG21	2:C:269:LYS:HD3	1.85	0.58
2:B:108:PRO:HB2	2:B:132:PRO:HA	1.85	0.58
2:D:40:ARG:HD3	2:D:85:CYS:HA	1.84	0.58
2:B:57:LEU:O	2:B:61:LYS:HG2	2.03	0.58
2:B:84:ASN:HD22	2:B:180:ASN:ND2	2.01	0.58
2:B:9:PRO:HD2	2:B:152:ILE:HD12	1.86	0.57
2:A:113:SER:O	2:A:149:GLY:HA2	2.02	0.57
2:B:36:VAL:HB	2:B:89:LEU:HB2	1.86	0.57
2:C:166:GLU:OE2	2:D:1:SER:HB3	2.05	0.57
2:B:40:ARG:HD3	2:B:85:CYS:HA	1.85	0.56
2:C:20:VAL:HG22	2:C:68:VAL:HG22	1.88	0.56
2:D:51:ASN:HB2	2:D:188:ARG:NE	2.19	0.56
1:F:0:GLN:HB3	2:D:142:ASN:HA	1.88	0.56
2:A:111:THR:HG22	2:A:129:ALA:HB2	1.86	0.56
2:C:20:VAL:HG12	2:C:42:VAL:HG21	1.87	0.55
2:D:222:ARG:HE	2:D:222:ARG:H	1.54	0.55
1:F:-1:LEU:HD12	2:D:41:HIS:CD2	2.41	0.55
2:A:44:CYS:HB3	2:A:48:ASP:HB2	1.88	0.55
1:F:-5:THR:HB	2:D:168:PRO:HB3	1.88	0.55
2:C:207:TRP:NE1	2:C:282:LEU:HD12	2.21	0.54
2:B:175:THR:HA	2:B:181:PHE:HA	1.90	0.54
2:A:58:LEU:HD11	2:A:80:HIS:CD2	2.40	0.54
2:B:243:THR:HG23	2:B:246:HIS:ND1	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:0:GLN:CB	2:D:142:ASN:HA	2.37	0.54
2:C:118:TYR:CE1	2:C:144:SER:HB3	2.43	0.53
2:B:64:HIS:CD2	2:B:65:ASN:ND2	2.77	0.53
2:C:102:LYS:HG2	2:C:104:VAL:HG23	1.91	0.53
2:C:201:THR:HG22	2:C:242:LEU:HD13	1.90	0.53
2:C:70:ALA:HB2	2:C:75:LEU:HD11	1.91	0.52
2:D:186:VAL:H	2:D:192:GLN:NE2	2.07	0.52
2:B:40:ARG:O	2:B:43:ILE:HG12	2.10	0.52
2:D:188:ARG:HB3	2:D:190:THR:HG22	1.91	0.52
2:D:43:ILE:HG22	2:D:66:PHE:HE1	1.75	0.52
2:D:271:LEU:HD11	2:D:281:ILE:HD11	1.91	0.52
2:B:14:GLU:HG2	2:B:122:PRO:HG2	1.92	0.52
2:B:85:CYS:HB2	2:B:175:THR:HG21	1.92	0.51
2:B:279:ARG:HH12	2:C:102:LYS:NZ	2.08	0.51
2:B:165:MET:HE1	2:B:185:PHE:HB3	1.93	0.51
2:D:276:MET:CE	2:D:281:ILE:HG13	2.40	0.51
1:E:-1:LEU:HD12	2:B:41:HIS:CD2	2.45	0.51
2:C:85:CYS:HB2	2:C:175:THR:HG21	1.93	0.51
2:C:10:SER:HB2	2:C:115:LEU:HD13	1.93	0.50
2:D:276:MET:HE3	2:D:281:ILE:HG13	1.92	0.50
2:D:164:HIS:CD2	2:D:175:THR:HG23	2.48	0.49
2:A:126:TYR:CE2	2:B:6:MET:HE2	2.46	0.49
2:C:2:GLY:HA3	2:D:139:SER:HA	1.95	0.49
2:B:206:ALA:HB2	2:B:293:PRO:HG3	1.95	0.49
2:B:222:ARG:HG2	2:C:107:GLN:HE21	1.77	0.49
2:C:21:THR:HB	2:C:67:LEU:HB2	1.94	0.49
2:A:245:ASP:O	2:A:249:ILE:HG13	2.13	0.49
2:C:100:LYS:HB2	2:C:156:CYS:HA	1.95	0.49
2:D:205:LEU:HD23	2:D:208:LEU:HD12	1.95	0.49
2:D:204:VAL:HG22	2:D:289:ASP:HB3	1.96	0.48
2:D:131:ARG:HG2	2:D:135:THR:O	2.14	0.48
2:D:222:ARG:H	2:D:222:ARG:NE	2.11	0.48
2:C:53:ASN:OD1	2:C:55:GLU:OE1	2.31	0.48
2:B:44:CYS:HB3	2:B:48:ASP:HB2	1.94	0.48
2:D:205:LEU:O	2:D:208:LEU:HB2	2.13	0.48
2:D:152:ILE:HG22	2:D:153:ASP:N	2.28	0.48
2:C:44:CYS:SG	2:C:54:TYR:CE1	3.07	0.47
2:C:132:PRO:HG2	2:C:198:THR:O	2.14	0.47
2:B:84:ASN:HD21	2:B:180:ASN:HD21	1.62	0.47
2:C:207:TRP:HZ3	2:C:287:LEU:HD23	1.79	0.47
2:B:108:PRO:CB	2:B:132:PRO:HA	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:106:ILE:HG12	2:C:160:CYS:HB2	1.97	0.47
2:A:95:ASN:HB3	2:A:98:THR:OG1	2.15	0.47
1:F:-5:THR:HB	2:D:168:PRO:CB	2.45	0.47
2:B:167:LEU:HB3	2:B:168:PRO:HD2	1.96	0.46
2:D:111:THR:HG23	2:D:292:THR:HG23	1.97	0.46
2:B:60:ARG:HG3	2:B:61:LYS:HE3	1.98	0.46
2:A:163:HIS:HE1	2:A:172:HIS:HB3	1.80	0.46
2:C:27:LEU:HD13	2:C:39:PRO:HD2	1.96	0.46
2:B:81:SER:HB3	2:B:88:LYS:HB2	1.96	0.46
2:D:138:GLY:O	2:D:172:HIS:HE1	1.98	0.46
2:C:3:PHE:HE1	2:C:300:CYS:HG	1.63	0.46
2:C:105:ARG:HH11	2:C:176:ASP:HB3	1.81	0.46
2:B:217:ARG:HB3	2:B:220:LEU:HD12	1.98	0.45
2:D:30:LEU:HD22	2:D:148:VAL:HG11	1.98	0.45
2:D:39:PRO:HG2	2:D:145:ALA:HB1	1.97	0.45
2:A:166:GLU:OE1	2:A:172:HIS:NE2	2.43	0.45
2:A:164:HIS:O	2:C:305:PHE:O	2.35	0.45
1:F:-1:LEU:CD1	2:D:41:HIS:CD2	2.99	0.45
2:C:3:PHE:CE1	2:C:300:CYS:SG	3.10	0.45
2:A:84:ASN:HB2	2:A:179:GLY:HA3	1.99	0.44
2:C:40:ARG:CD	2:C:85:CYS:HA	2.44	0.44
2:B:210:ALA:HB2	2:B:296:VAL:HG13	1.98	0.44
2:B:270:GLU:OE2	2:B:273:GLN:OE1	2.35	0.44
2:B:163:HIS:HE1	2:B:172:HIS:HB3	1.81	0.44
2:A:163:HIS:CE1	2:A:172:HIS:HB3	2.53	0.44
2:A:91:VAL:HG23	2:A:93:THR:H	1.83	0.44
2:D:17:MET:HG3	2:D:117:CYS:SG	2.58	0.44
2:D:51:ASN:HB2	2:D:188:ARG:HH21	1.82	0.44
2:D:269:LYS:HA	2:D:272:LEU:HD12	1.99	0.44
2:D:249:ILE:HG22	2:D:293:PRO:HG2	1.98	0.44
2:D:165:MET:HE2	2:D:173:ALA:HB3	2.00	0.43
2:D:53:ASN:HB3	2:D:56:ASP:HB2	2.00	0.43
2:D:212:VAL:HG21	2:D:264:MET:HE1	2.00	0.43
2:B:198:THR:HA	2:B:238:ASN:OD1	2.18	0.43
2:D:163:HIS:HE1	2:D:172:HIS:HB3	1.83	0.43
2:B:249:ILE:HG22	2:B:293:PRO:HG2	1.99	0.43
2:D:12:LYS:NZ	2:D:99:PRO:HB3	2.33	0.43
2:C:185:PHE:HD1	2:C:192:GLN:HG2	1.84	0.43
2:A:164:HIS:CD2	2:A:175:THR:HG23	2.54	0.43
2:B:52:PRO:HD3	2:B:188:ARG:HG3	2.01	0.43
2:D:294:PHE:O	2:D:298:ARG:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:8:PHE:HB3	2:A:152:ILE:HD12	2.00	0.42
2:C:234:ALA:O	2:C:239:TYR:HB2	2.19	0.42
2:A:262:LEU:HA	2:A:265:CYS:HB2	2.01	0.42
2:B:84:ASN:ND2	2:B:180:ASN:HD22	2.16	0.42
2:B:95:ASN:HB3	2:B:98:THR:OG1	2.18	0.42
2:C:117:CYS:HA	2:C:122:PRO:HA	2.00	0.42
2:D:54:TYR:HA	2:D:57:LEU:HD12	2.00	0.42
2:B:31:TRP:CD2	2:B:95:ASN:HB2	2.54	0.42
2:B:279:ARG:NH1	2:C:102:LYS:NZ	2.67	0.42
2:C:101:TYR:O	2:C:156:CYS:SG	2.78	0.42
2:C:111:THR:HG22	2:C:129:ALA:HB2	2.01	0.42
2:D:10:SER:O	2:D:14:GLU:HG3	2.20	0.42
2:C:140:PHE:HB2	2:C:172:HIS:CD2	2.55	0.42
2:A:109:GLY:HA2	2:A:200:ILE:HD13	2.01	0.41
2:C:294:PHE:O	2:C:298:ARG:HB2	2.19	0.41
2:D:188:ARG:HH11	2:D:190:THR:HB	1.85	0.41
2:B:165:MET:CE	2:B:185:PHE:HB3	2.50	0.41
2:B:209:TYR:HB3	2:B:253:LEU:HD22	2.03	0.41
2:B:140:PHE:HB3	2:B:144:SER:OG	2.20	0.41
2:D:299:GLN:HE21	2:D:299:GLN:HB2	1.64	0.41
2:D:27:LEU:HD13	2:D:39:PRO:HD2	2.02	0.41
2:A:167:LEU:HB3	2:A:168:PRO:HD2	2.03	0.41
2:C:52:PRO:HD3	2:C:188:ARG:HG3	2.02	0.41
2:D:186:VAL:H	2:D:192:GLN:HE22	1.67	0.41
2:A:27:LEU:HB2	2:A:145:ALA:O	2.21	0.41
1:E:-1:LEU:HD23	2:B:189:GLN:HG3	2.03	0.41
2:A:216:ASP:O	2:A:217:ARG:HG3	2.21	0.41
2:D:161:TYR:CE1	2:D:174:GLY:HA3	2.55	0.41
2:B:233:VAL:HG21	2:B:269:LYS:HD3	2.03	0.40
2:D:47:GLU:HG2	2:D:50:LEU:HD12	2.03	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	E	4/6 (67%)	4 (100%)	0	0	100	100
1	F	4/6 (67%)	3 (75%)	1 (25%)	0	100	100
2	A	304/306 (99%)	280 (92%)	19 (6%)	5 (2%)	7	31
2	B	304/306 (99%)	274 (90%)	26 (9%)	4 (1%)	9	36
2	C	304/306 (99%)	265 (87%)	36 (12%)	3 (1%)	12	42
2	D	303/306 (99%)	264 (87%)	38 (12%)	1 (0%)	36	67
All	All	1223/1236 (99%)	1090 (89%)	120 (10%)	13 (1%)	11	40

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	154	TYR
2	B	47	GLU
2	C	154	TYR
2	B	63	ASN
2	A	155	ASP
2	B	43	ILE
2	C	283	GLY
2	A	50	LEU
2	C	134	PHE
2	D	240	GLU
2	A	2	GLY
2	A	11	GLY
2	B	52	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	E	5/5 (100%)	3 (60%)	2 (40%)	0	0
1	F	5/5 (100%)	4 (80%)	1 (20%)	1	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	A	262/262 (100%)	239 (91%)	23 (9%)	9	33
2	B	262/262 (100%)	243 (93%)	19 (7%)	13	40
2	C	262/262 (100%)	227 (87%)	35 (13%)	4	16
2	D	261/262 (100%)	227 (87%)	34 (13%)	4	17
All	All	1057/1058 (100%)	943 (89%)	114 (11%)	6	24

All (114) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	-5	THR
1	E	0	GLN
2	A	1	SER
2	A	4	ARG
2	A	5	LYS
2	A	25	THR
2	A	44	CYS
2	A	45	THR
2	A	55	GLU
2	A	64	HIS
2	A	67	LEU
2	A	87	LEU
2	A	102	LYS
2	A	106	ILE
2	A	155	ASP
2	A	169	THR
2	A	188	ARG
2	A	190	THR
2	A	192	GLN
2	A	197	ASP
2	A	227	LEU
2	A	228	ASN
2	A	288	GLU
2	A	297	VAL
2	A	299	GLN
2	B	35	VAL
2	B	47	GLU
2	B	49	MET
2	B	55	GLU
2	B	58	LEU
2	B	67	LEU
2	B	69	GLN

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Mol	Chain	Res	Type
2	B	73	VAL
2	B	74	GLN
2	B	87	LEU
2	B	107	GLN
2	B	128	CYS
2	B	154	TYR
2	B	158	SER
2	B	216	ASP
2	B	224	THR
2	B	227	LEU
2	B	228	ASN
2	B	232	LEU
1	F	0	GLN
2	C	1	SER
2	C	4	ARG
2	C	16	CYS
2	C	19	GLN
2	C	22	CYS
2	C	35	VAL
2	C	48	ASP
2	C	58	LEU
2	C	60	ARG
2	C	62	SER
2	C	64	HIS
2	C	73	VAL
2	C	80	HIS
2	C	84	ASN
2	C	87	LEU
2	C	89	LEU
2	C	90	LYS
2	C	106	ILE
2	C	110	GLN
2	C	114	VAL
2	C	128	CYS
2	C	162	MET
2	C	177	LEU
2	C	192	GLN
2	C	217	ARG
2	C	225	THR
2	C	248	ASP
2	C	269	LYS
2	C	280	THR

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Mol	Chain	Res	Type
2	C	286	LEU
2	C	288	GLU
2	C	289	ASP
2	C	291	PHE
2	C	292	THR
2	C	297	VAL
2	D	6	MET
2	D	10	SER
2	D	12	LYS
2	D	35	VAL
2	D	48	ASP
2	D	50	LEU
2	D	56	ASP
2	D	73	VAL
2	D	78	ILE
2	D	87	LEU
2	D	104	VAL
2	D	107	GLN
2	D	121	SER
2	D	128	CYS
2	D	142	ASN
2	D	158	SER
2	D	165	MET
2	D	186	VAL
2	D	188	ARG
2	D	200	ILE
2	D	204	VAL
2	D	222	ARG
2	D	224	THR
2	D	225	THR
2	D	227	LEU
2	D	233	VAL
2	D	242	LEU
2	D	243	THR
2	D	245	ASP
2	D	263	ASP
2	D	271	LEU
2	D	284	SER
2	D	286	LEU
2	D	304	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	51	ASN
2	A	53	ASN
2	A	256	GLN
2	A	274	ASN
2	B	41	HIS
2	B	64	HIS
2	B	65	ASN
2	B	84	ASN
2	B	110	GLN
2	B	180	ASN
2	B	214	ASN
2	B	228	ASN
2	B	256	GLN
2	B	274	ASN
1	F	0	GLN
2	C	64	HIS
2	C	80	HIS
2	C	151	ASN
2	C	214	ASN
2	D	41	HIS
2	D	107	GLN
2	D	192	GLN
2	D	299	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 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	E	6/6 (100%)	0.67	0 100 100	81, 82, 83, 83	0
1	F	6/6 (100%)	0.91	1 (16%) 4 3	74, 78, 84, 85	0
2	A	306/306 (100%)	-0.08	2 (0%) 84 70	30, 48, 67, 79	0
2	B	306/306 (100%)	0.18	3 (0%) 79 63	33, 59, 88, 98	0
2	C	306/306 (100%)	0.60	7 (2%) 61 42	45, 71, 92, 103	0
2	D	305/306 (99%)	0.27	5 (1%) 70 51	41, 65, 94, 108	0
All	All	1235/1236 (99%)	0.24	18 (1%) 72 53	30, 61, 90, 108	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	155	ASP	3.1
2	D	188	ARG	3.1
2	C	294	PHE	2.9
2	C	287	LEU	2.8
2	B	275	GLY	2.4
2	C	150	PHE	2.4
2	C	206	ALA	2.4
2	D	201	THR	2.4
2	C	50	LEU	2.4
2	C	58	LEU	2.3
2	D	230	PHE	2.3
1	F	-1	LEU	2.2
2	D	194	ALA	2.2
2	D	154	TYR	2.2
2	B	66	PHE	2.1
2	B	256	GLN	2.1
2	A	303	VAL	2.1
2	C	280	THR	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.