



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:52 AM UTC

PDB ID : 7F7H / pdb_00007f7h
Title : SARS-CoV-2 S protein RBD in complex with A8-1 Fab
Authors : Dou, Y.; Wang, X.; Wang, K.; Liu, P.; Lu, B.
Deposited on : 2021-06-29
Resolution : 3.19 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

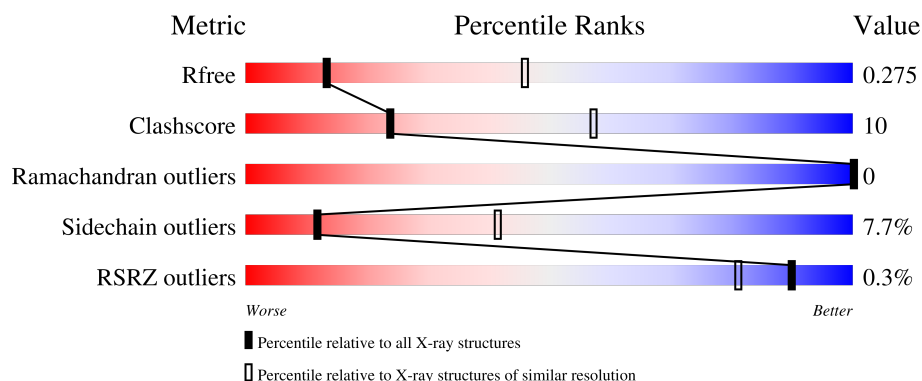
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	218	 69% 27% . .
1	C	218	 74% 23% .
2	B	212	 75% 23% .
2	D	212	 73% 23% . .
3	E	182	 75% 21% .

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Mol	Chain	Length	Quality of chain
3	F	182	2% 67% 26% 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9218 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 Heavy chain of A8-1 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	0	1
			1572	993	261	311	7			
1	C	218	Total	C	N	O	S	0	0	0
			1615	1018	270	320	7			

- Molecule 2 is a protein called Light chain of A8-1 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	0	0
			1616	1014	272	326	4			
2	D	210	Total	C	N	O	S	0	0	0
			1603	1007	270	322	4			

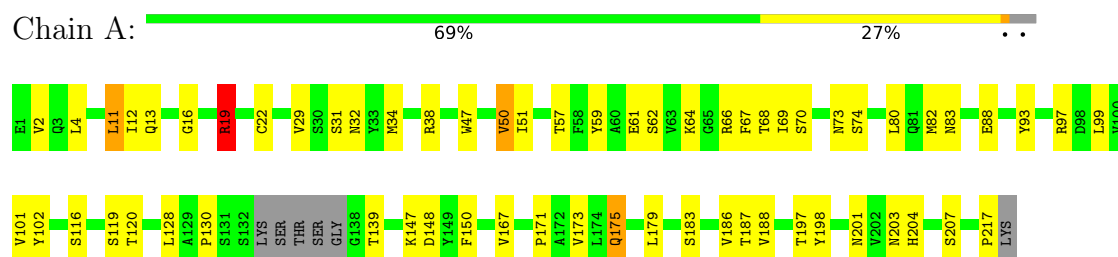
- Molecule 3 is a protein called Spike glycoprotein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	182	Total	C	N	O	S	0	0	0
			1453	932	242	272	7			
3	F	170	Total	C	N	O	S	0	0	0
			1359	872	227	254	6			

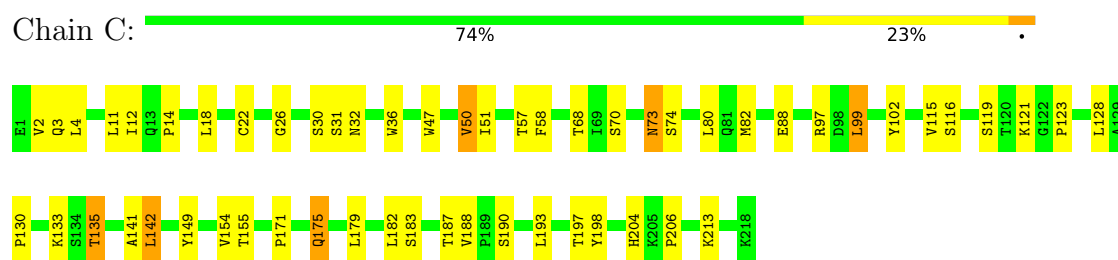
3 Residue-property plots [i](#)

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.

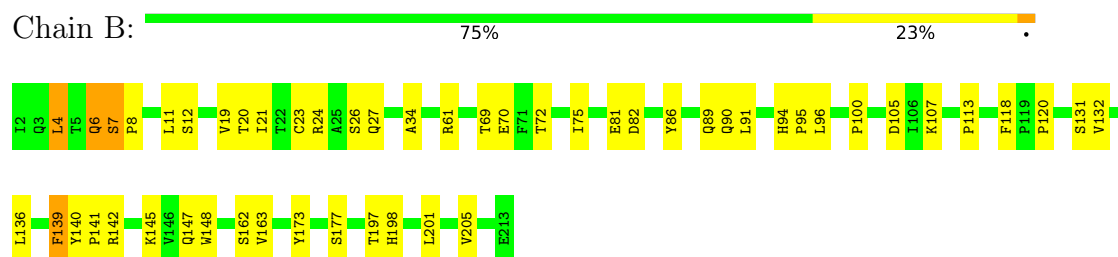
- Molecule 1: Heavy chain of A8-1 Fab



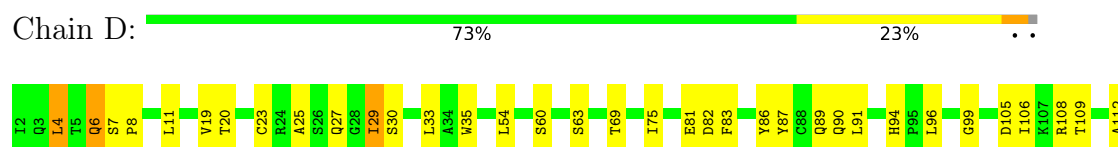
- Molecule 1: Heavy chain of A8-1 Fab

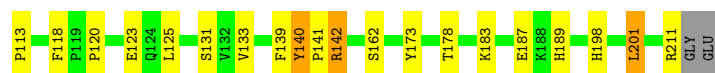


- Molecule 2: Light chain of A8-1 Fab

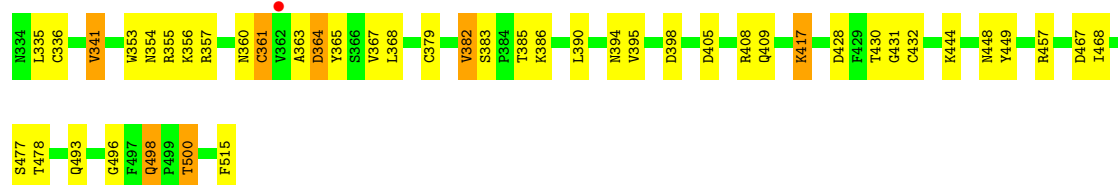
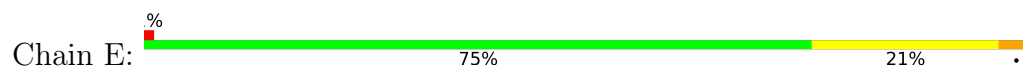


- Molecule 2: Light chain of A8-1 Fab

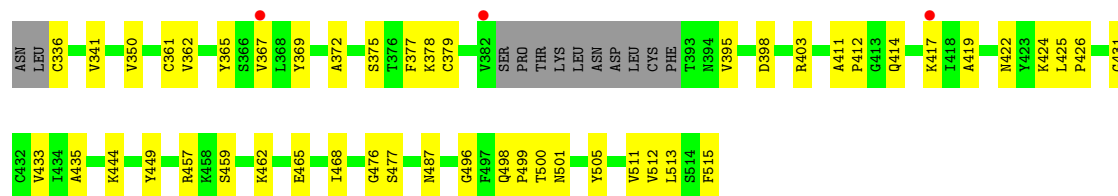




• Molecule 3: Spike glycoprotein S1



• Molecule 3: Spike glycoprotein S1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.77Å 75.66Å 181.54Å 90.00° 92.26° 90.00°	Depositor
Resolution (Å)	36.28 – 3.19 36.28 – 3.19	Depositor EDS
% Data completeness (in resolution range)	99.3 (36.28-3.19) 99.2 (36.28-3.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 3.18Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.205 , 0.275 0.205 , 0.275	Depositor DCC
R_{free} test set	1195 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	68.9	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 36.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.086 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9218	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% 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	0.48	0/1606	0.67	1/2189 (0.0%)
1	C	0.43	0/1650	0.65	0/2248
2	B	0.48	0/1651	0.77	4/2243 (0.2%)
2	D	0.45	0/1638	0.71	1/2226 (0.0%)
3	E	0.45	0/1494	0.65	0/2031
3	F	0.43	0/1397	0.63	0/1897
All	All	0.45	0/9436	0.68	6/12834 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	139	PHE	CA-C-N	-5.67	107.96	121.80
2	B	139	PHE	C-N-CA	-5.67	107.96	121.80
2	D	140	TYR	CA-CB-CG	-5.33	104.32	113.90
2	B	6	GLN	CA-C-N	5.22	125.62	119.83
2	B	6	GLN	C-N-CA	5.22	125.62	119.83
1	A	19	ARG	CG-CD-NE	5.15	123.34	112.00

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	1572	0	1553	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1615	0	1602	34	0
2	B	1616	0	1580	33	0
2	D	1603	0	1571	36	0
3	E	1453	0	1374	28	0
3	F	1359	0	1278	29	0
All	All	9218	0	8958	183	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 10.

All (183) 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)
3:F:336:CYS:N	3:F:365:TYR:HH	1.34	1.21
3:F:336:CYS:N	3:F:365:TYR:OH	2.13	0.80
2:B:141:PRO:HD2	2:B:198:HIS:CE1	2.17	0.79
1:A:204:HIS:HD1	1:A:207:SER:HG	1.34	0.73
1:C:141:ALA:HB2	1:C:187:THR:HG22	1.71	0.71
2:B:141:PRO:HD2	2:B:198:HIS:HE1	1.55	0.71
2:D:141:PRO:HD2	2:D:198:HIS:CE1	2.27	0.70
1:C:213:LYS:NZ	2:D:123:GLU:OE1	2.24	0.70
1:A:19:ARG:HB3	1:A:19:ARG:HH11	1.58	0.69
2:B:145:LYS:HB3	2:B:197:THR:HB	1.75	0.68
2:D:120:PRO:HB3	2:D:131:SER:H	1.58	0.68
2:B:142:ARG:HH22	2:B:163:VAL:HG21	1.57	0.68
1:C:47:TRP:HE1	1:C:50:VAL:HG12	1.57	0.68
1:A:12:ILE:HD11	1:A:16:GLY:HA3	1.75	0.67
3:E:355:ARG:NH2	3:E:398:ASP:OD1	2.26	0.67
2:D:140:TYR:HB3	2:D:141:PRO:HD3	1.77	0.67
1:A:171:PRO:HD2	2:B:162:SER:HB3	1.76	0.66
2:B:140:TYR:O	2:B:142:ARG:N	2.28	0.66
2:D:6:GLN:NE2	2:D:86:TYR:O	2.31	0.63
2:B:142:ARG:O	2:B:142:ARG:HD3	1.98	0.63
3:E:457:ARG:NH1	3:E:467:ASP:OD2	2.32	0.62
3:F:431:GLY:HA2	3:F:515:PHE:CE2	2.35	0.62
2:B:142:ARG:NH2	2:B:163:VAL:HG11	2.15	0.62
2:B:120:PRO:HB3	2:B:131:SER:H	1.65	0.61
1:A:47:TRP:HB2	2:B:96:LEU:HB2	1.82	0.61
1:C:73:ASN:OD1	1:C:73:ASN:N	2.27	0.60
1:A:167:VAL:HG22	1:A:186:VAL:HB	1.84	0.59
2:B:113:PRO:HB3	2:B:139:PHE:HB3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:409:GLN:HE22	3:E:417:LYS:HD2	1.67	0.59
3:F:431:GLY:HA3	3:F:513:LEU:O	2.02	0.58
2:D:6:GLN:HE22	2:D:87:TYR:HA	1.69	0.58
2:B:61:ARG:NH2	2:B:81:GLU:OE2	2.37	0.57
1:C:135:THR:HB	1:C:190:SER:HB2	1.85	0.57
1:C:99:LEU:HB2	1:C:102:TYR:HB2	1.87	0.57
2:B:201:LEU:HD13	2:B:205:VAL:HG23	1.88	0.56
3:F:367:VAL:HG23	3:F:369:TYR:CD1	2.40	0.56
2:B:140:TYR:HB3	2:B:141:PRO:HD3	1.87	0.56
3:E:449:TYR:CE1	3:E:496:GLY:HA2	2.41	0.56
3:F:457:ARG:NH1	3:F:459:SER:O	2.37	0.56
3:E:449:TYR:HE1	3:E:496:GLY:HA2	1.70	0.56
1:A:204:HIS:ND1	1:A:207:SER:OG	2.26	0.55
1:A:29:VAL:HG12	1:A:34:MET:HE2	1.87	0.55
1:C:36:TRP:NE1	1:C:80:LEU:HB2	2.22	0.55
1:A:13:GLN:H	1:A:13:GLN:CD	2.15	0.55
1:C:47:TRP:HB2	2:D:96:LEU:HB2	1.89	0.55
3:F:403:ARG:HD2	3:F:505:TYR:HA	1.89	0.55
2:D:83:PHE:CZ	2:D:106:ILE:HG12	2.43	0.54
3:F:336:CYS:HB2	3:F:365:TYR:CE2	2.43	0.54
1:A:128:LEU:HB3	2:B:118:PHE:CD2	2.43	0.54
2:B:105:ASP:OD1	2:B:173:TYR:OH	2.25	0.54
1:A:51:ILE:HG13	1:A:57:THR:HG22	1.90	0.53
2:D:30:SER:HB3	3:F:501:ASN:HD21	1.72	0.53
1:C:97:ARG:CZ	1:C:99:LEU:HD21	2.38	0.53
1:C:154:VAL:HG12	1:C:204:HIS:HB2	1.90	0.53
1:C:30:SER:OG	1:C:73:ASN:HB3	2.09	0.53
1:A:175:GLN:HG2	1:A:179:LEU:O	2.07	0.53
1:C:4:LEU:HB3	1:C:22:CYS:SG	2.50	0.52
1:C:12:ILE:HG23	1:C:115:VAL:HG22	1.90	0.52
2:D:23:CYS:HB2	2:D:35:TRP:CH2	2.44	0.52
2:D:141:PRO:HD2	2:D:198:HIS:HE1	1.75	0.52
1:A:97:ARG:CZ	1:A:99:LEU:HD21	2.41	0.51
3:E:379:CYS:HA	3:E:432:CYS:HA	1.93	0.51
3:F:417:LYS:HD3	3:F:417:LYS:H	1.75	0.51
2:D:105:ASP:OD1	2:D:173:TYR:OH	2.28	0.51
3:F:449:TYR:HE1	3:F:496:GLY:HA2	1.75	0.51
1:A:188:VAL:HG11	1:A:198:TYR:CE1	2.46	0.50
2:D:112:ALA:HB1	2:D:201:LEU:HD12	1.92	0.50
2:D:19:VAL:HG22	2:D:75:ILE:HB	1.93	0.50
2:B:7:SER:OG	2:B:8:PRO:HD3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:19:VAL:HG22	2:B:75:ILE:HB	1.94	0.50
1:A:19:ARG:HH11	1:A:19:ARG:CB	2.25	0.50
3:E:354:ASN:O	3:E:398:ASP:HA	2.12	0.50
2:D:6:GLN:OE1	2:D:99:GLY:HA3	2.12	0.50
3:F:375:SER:HB3	3:F:433:VAL:HB	1.94	0.50
3:E:341:VAL:HG12	3:E:356:LYS:HD3	1.93	0.49
2:D:54:LEU:HD21	2:D:60:SER:HA	1.95	0.48
1:C:171:PRO:HD2	2:D:162:SER:HB3	1.96	0.48
3:F:425:LEU:HD21	3:F:512:VAL:HG11	1.94	0.48
1:A:4:LEU:HB3	1:A:22:CYS:SG	2.53	0.48
2:D:113:PRO:HD2	2:D:201:LEU:CD1	2.44	0.48
2:B:105:ASP:HB3	2:B:140:TYR:OH	2.14	0.48
1:C:14:PRO:HD3	1:C:116:SER:O	2.14	0.47
2:D:82:ASP:O	2:D:86:TYR:OH	2.23	0.47
3:E:368:LEU:HA	3:E:368:LEU:HD23	1.75	0.47
3:F:476:GLY:HA3	3:F:487:ASN:HD22	1.80	0.47
3:E:353:TRP:H	3:E:353:TRP:CD1	2.31	0.47
3:E:408:ARG:HE	3:E:408:ARG:HB3	1.50	0.47
2:B:142:ARG:NH2	2:B:163:VAL:HG21	2.27	0.47
3:F:362:VAL:HB	3:F:365:TYR:OH	2.15	0.47
1:A:88:GLU:N	1:A:88:GLU:OE1	2.48	0.47
2:B:82:ASP:O	2:B:86:TYR:OH	2.24	0.47
3:E:444:LYS:HG2	3:E:448:ASN:HB2	1.97	0.46
1:A:99:LEU:HB2	1:A:102:TYR:HB2	1.97	0.46
1:C:32:ASN:OD1	1:C:97:ARG:HD3	2.15	0.46
1:C:128:LEU:HB3	2:D:118:PHE:CD2	2.50	0.46
1:C:175:GLN:NE2	1:C:179:LEU:O	2.42	0.46
2:D:4:LEU:HG	2:D:23:CYS:SG	2.55	0.46
3:F:411:ALA:HB3	3:F:414:GLN:HG3	1.97	0.46
3:E:386:LYS:HB3	3:E:386:LYS:HE2	1.53	0.46
2:D:113:PRO:HB3	2:D:139:PHE:HB3	1.98	0.46
1:C:130:PRO:HB3	1:C:142:LEU:HB3	1.97	0.46
1:A:13:GLN:HA	1:A:116:SER:O	2.15	0.46
1:C:18:LEU:HB3	1:C:82:MET:HE3	1.97	0.46
1:A:62:SER:O	1:A:66:ARG:NH2	2.49	0.46
2:D:7:SER:HB3	2:D:8:PRO:HD3	1.97	0.45
1:A:11:LEU:HD11	1:A:150:PHE:CE2	2.51	0.45
3:E:428:ASP:O	3:E:430:THR:HG23	2.16	0.45
2:D:183:LYS:O	2:D:187:GLU:HG2	2.16	0.45
2:D:189:HIS:O	2:D:211:ARG:NH1	2.39	0.45
1:A:67:PHE:CD1	1:A:82:MET:HA	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:PRO:HG2	1:A:217:PRO:HB3	1.98	0.45
3:F:367:VAL:HG23	3:F:369:TYR:HD1	1.80	0.45
3:F:398:ASP:O	3:F:511:VAL:HA	2.16	0.45
3:E:335:LEU:HG	3:E:336:CYS:H	1.81	0.44
2:D:133:VAL:HG22	2:D:178:THR:HG23	1.98	0.44
3:F:412:PRO:HB3	3:F:426:PRO:O	2.17	0.44
1:A:120:THR:HA	1:A:150:PHE:O	2.16	0.44
1:C:190:SER:HA	1:C:193:LEU:HG	2.00	0.44
2:B:89:GLN:HG2	2:B:90:GLN:N	2.32	0.43
3:E:405:ASP:N	3:E:405:ASP:OD1	2.48	0.43
1:A:66:ARG:HB3	1:A:83:ASN:O	2.17	0.43
2:B:24:ARG:NH1	2:B:70:GLU:OE1	2.52	0.43
3:E:336:CYS:SG	3:E:360:ASN:HB2	2.58	0.43
1:C:133:LYS:HD3	1:C:133:LYS:HA	1.77	0.43
2:B:94:HIS:CD2	2:B:95:PRO:HD3	2.54	0.43
3:E:382:VAL:HG22	3:E:383:SER:O	2.17	0.43
2:B:142:ARG:HG3	2:B:173:TYR:CE2	2.54	0.43
3:E:498:GLN:HB3	3:E:500:THR:HG23	2.01	0.43
2:D:125:LEU:O	2:D:183:LYS:HD2	2.18	0.43
1:C:88:GLU:N	1:C:88:GLU:OE1	2.52	0.43
2:D:89:GLN:NE2	2:D:90:GLN:O	2.47	0.43
1:A:101:VAL:HG11	3:E:493:GLN:HG3	2.01	0.43
2:D:142:ARG:O	2:D:142:ARG:HD2	2.19	0.43
3:F:417:LYS:H	3:F:417:LYS:CD	2.31	0.43
1:A:148:ASP:HB3	1:A:179:LEU:HD13	2.01	0.43
1:A:80:LEU:HA	1:A:80:LEU:HD12	1.70	0.42
3:E:363:ALA:HB3	3:E:365:TYR:CE1	2.53	0.42
1:A:59:TYR:OH	1:A:69:ILE:HG22	2.19	0.42
2:D:108:ARG:HD3	2:D:109:THR:O	2.19	0.42
3:F:444:LYS:O	3:F:499:PRO:HD3	2.18	0.42
2:B:94:HIS:HB2	3:E:417:LYS:NZ	2.35	0.42
2:B:34:ALA:HB2	2:B:91:LEU:HD21	2.01	0.42
1:C:133:LYS:HA	1:C:135:THR:HG23	2.01	0.42
2:D:112:ALA:HB1	2:D:201:LEU:CD1	2.49	0.42
2:B:6:GLN:HG3	2:B:100:PRO:HD2	2.02	0.42
2:B:4:LEU:HG	2:B:23:CYS:SG	2.59	0.42
1:A:32:ASN:OD1	1:A:97:ARG:HD3	2.20	0.42
1:A:188:VAL:HG11	1:A:198:TYR:CZ	2.55	0.42
2:D:29:ILE:O	2:D:30:SER:C	2.62	0.42
3:E:360:ASN:HB3	3:E:361:CYS:H	1.56	0.42
3:F:372:ALA:HA	3:F:435:ALA:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:GLY:O	3:F:487:ASN:ND2	2.50	0.41
2:B:148:TRP:NE1	2:B:177:SER:OG	2.44	0.41
1:C:182:LEU:C	1:C:182:LEU:HD12	2.46	0.41
3:F:419:ALA:O	3:F:424:LYS:HE2	2.20	0.41
1:C:188:VAL:HG11	1:C:198:TYR:CZ	2.55	0.41
3:F:377:PHE:CZ	3:F:412:PRO:HD2	2.55	0.41
3:E:409:GLN:NE2	3:E:417:LYS:HD2	2.32	0.41
3:F:403:ARG:HH11	3:F:403:ARG:HD3	1.73	0.41
2:B:12:SER:HB3	2:B:107:LYS:HG3	2.02	0.41
3:F:350:VAL:HG22	3:F:422:ASN:HB3	2.02	0.41
3:E:364:ASP:HB3	3:E:367:VAL:HG21	2.02	0.41
1:C:36:TRP:CE2	1:C:80:LEU:HB2	2.55	0.41
1:C:133:LYS:C	1:C:135:THR:H	2.29	0.41
1:C:193:LEU:HD23	1:C:193:LEU:HA	1.85	0.41
2:D:4:LEU:HD12	2:D:25:ALA:HA	2.02	0.41
2:D:83:PHE:CE1	2:D:106:ILE:HG12	2.55	0.41
1:C:123:PRO:HB3	1:C:149:TYR:HB3	2.02	0.41
2:D:4:LEU:HD21	2:D:33:LEU:HD11	2.02	0.41
3:F:378:LYS:HD3	3:F:378:LYS:HA	1.63	0.41
3:F:462:LYS:N	3:F:465:GLU:OE1	2.50	0.41
1:A:47:TRP:HE1	1:A:50:VAL:HG12	1.86	0.40
3:E:444:LYS:HB3	3:E:444:LYS:HE2	1.90	0.40
1:C:58:PHE:CG	2:D:94:HIS:HD2	2.39	0.40
1:C:204:HIS:CD2	1:C:206:PRO:HD2	2.56	0.40
1:C:51:ILE:HG13	1:C:57:THR:HG22	2.03	0.40
1:A:38:ARG:NH1	1:A:93:TYR:OH	2.51	0.40
2:B:21:ILE:O	2:B:72:THR:HA	2.22	0.40
1:A:19:ARG:HB3	1:A:19:ARG:NH1	2.30	0.40
1:A:61:GLU:HA	1:A:64:LYS:HE3	2.03	0.40
2:B:132:VAL:HG12	2:B:148:TRP:CH2	2.57	0.40
3:E:357:ARG:NH1	3:E:394:ASN:HD21	2.19	0.40
3:E:431:GLY:HA2	3:E:515:PHE:CE2	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	208/218 (95%)	199 (96%)	9 (4%)	0	100	100
1	C	216/218 (99%)	207 (96%)	9 (4%)	0	100	100
2	B	210/212 (99%)	198 (94%)	12 (6%)	0	100	100
2	D	208/212 (98%)	194 (93%)	14 (7%)	0	100	100
3	E	180/182 (99%)	163 (91%)	17 (9%)	0	100	100
3	F	166/182 (91%)	148 (89%)	18 (11%)	0	100	100
All	All	1188/1224 (97%)	1109 (93%)	79 (7%)	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	178/183 (97%)	159 (89%)	19 (11%)	6	27
1	C	183/183 (100%)	165 (90%)	18 (10%)	7	31
2	B	184/184 (100%)	175 (95%)	9 (5%)	22	55
2	D	183/184 (100%)	171 (93%)	12 (7%)	15	47
3	E	158/158 (100%)	145 (92%)	13 (8%)	10	39
3	F	146/158 (92%)	138 (94%)	8 (6%)	19	52
All	All	1032/1050 (98%)	953 (92%)	79 (8%)	12	41

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	11	LEU
1	A	19	ARG

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Mol	Chain	Res	Type
1	A	31	SER
1	A	50	VAL
1	A	68	THR
1	A	70	SER
1	A	73	ASN
1	A	74	SER
1	A	119	SER
1	A	139	THR
1	A	147	LYS
1	A	173	VAL
1	A	175	GLN
1	A	183	SER
1	A	187	THR
1	A	197	THR
1	A	201	ASN
1	A	203	ASN
2	B	4	LEU
2	B	7	SER
2	B	11	LEU
2	B	20	THR
2	B	26	SER
2	B	27	GLN
2	B	69	THR
2	B	136	LEU
2	B	147	GLN
3	E	341	VAL
3	E	361	CYS
3	E	364	ASP
3	E	382	VAL
3	E	385	THR
3	E	390	LEU
3	E	395	VAL
3	E	417	LYS
3	E	468	ILE
3	E	477	SER
3	E	478	THR
3	E	498	GLN
3	E	500	THR
1	C	2	VAL
1	C	3	GLN
1	C	11	LEU
1	C	31	SER

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Mol	Chain	Res	Type
1	C	50	VAL
1	C	68	THR
1	C	70	SER
1	C	73	ASN
1	C	74	SER
1	C	99	LEU
1	C	119	SER
1	C	121	LYS
1	C	135	THR
1	C	142	LEU
1	C	155	THR
1	C	175	GLN
1	C	183	SER
1	C	197	THR
2	D	4	LEU
2	D	6	GLN
2	D	11	LEU
2	D	20	THR
2	D	27	GLN
2	D	29	ILE
2	D	63	SER
2	D	69	THR
2	D	81	GLU
2	D	91	LEU
2	D	142	ARG
2	D	201	LEU
3	F	341	VAL
3	F	361	CYS
3	F	379	CYS
3	F	395	VAL
3	F	468	ILE
3	F	477	SER
3	F	498	GLN
3	F	500	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	27	GLN
2	B	94	HIS
2	B	147	GLN
2	B	198	HIS

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Mol	Chain	Res	Type
3	E	360	ASN
1	C	83	ASN
1	C	203	ASN
2	D	55	GLN
2	D	79	GLN
2	D	94	HIS
3	F	487	ASN
3	F	501	ASN

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	212/218 (97%)	-0.37	0	100	100	42, 54, 65, 77	0
1	C	218/218 (100%)	-0.31	0	100	100	48, 56, 75, 105	0
2	B	212/212 (100%)	-0.28	0	100	100	42, 53, 67, 83	0
2	D	210/212 (99%)	-0.21	0	100	100	49, 59, 71, 77	0
3	E	182/182 (100%)	-0.00	1 (0%)	87	76	44, 65, 104, 119	0
3	F	170/182 (93%)	0.01	3 (1%)	67	48	54, 71, 111, 124	0
All	All	1204/1224 (98%)	-0.20	4 (0%)	90	81	42, 58, 94, 124	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	417	LYS	2.6
3	F	382	VAL	2.2
3	E	362	VAL	2.2
3	F	367	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 ⓘ

There are no such residues in this entry.