



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:02 AM UTC

PDB ID : 6PXG / pdb_00006pxg
Title : Crystal Structure of MERS-CoV neutralizing antibody G2 Fab
Authors : Wang, N.; McLellan, J.S.
Deposited on : 2019-07-26
Resolution : 2.10 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

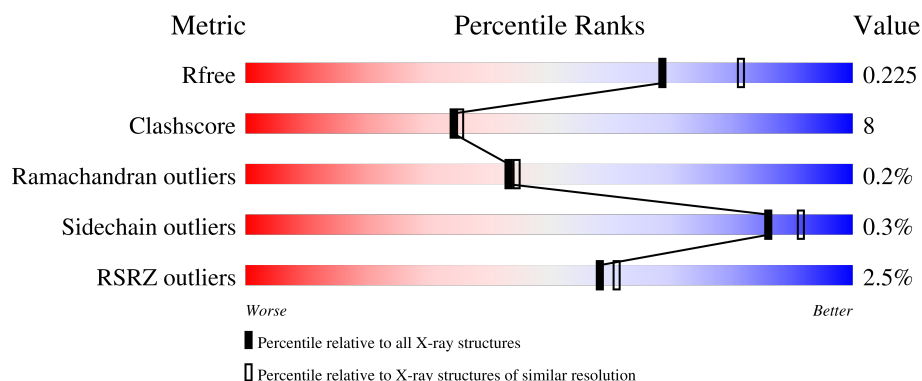
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	229	0% 81% 11% 8%
1	C	229	83% 9% 8%
1	E	229	6% 75% 17% 8%
1	H	229	0% 80% 12% 8%
2	B	218	2% 87% 12% •

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Mol	Chain	Length	Quality of chain
2	D	218	3% 79% 19% ••
2	F	218	2% 81% 18% •
2	L	218	2% 87% 12% •

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13415 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 G2 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	210	Total	C	N	O	S	0	0	0
			1570	998	250	315	7			
1	C	211	Total	C	N	O	S	0	0	0
			1579	1003	252	317	7			
1	E	210	Total	C	N	O	S	0	0	0
			1570	998	250	315	7			
1	H	210	Total	C	N	O	S	0	0	0
			1570	998	250	315	7			

- Molecule 2 is a protein called G2 Fab Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	216	Total	C	N	O	S	0	0	0
			1659	1030	283	339	7			
2	D	216	Total	C	N	O	S	0	0	0
			1659	1030	283	339	7			
2	F	216	Total	C	N	O	S	0	0	0
			1659	1030	283	339	7			
2	L	216	Total	C	N	O	S	0	0	0
			1659	1030	283	339	7			

- Molecule 3 is water.

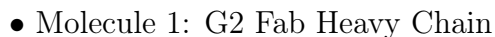
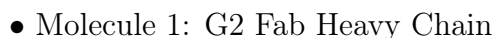
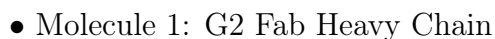
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total	O	0	0
			92	92		
3	B	80	Total	O	0	0
			80	80		
3	C	89	Total	O	0	0
			89	89		
3	D	48	Total	O	0	0
			48	48		

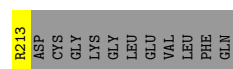
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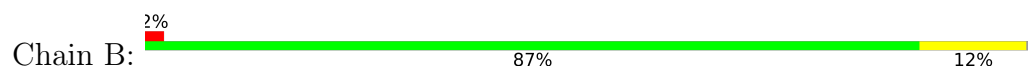
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	25	Total 25	O 25	0	0
3	F	42	Total 42	O 42	0	0
3	H	65	Total 65	O 65	0	0
3	L	49	Total 49	O 49	0	0

- Molecule 1: G2 Fab Heavy Chain

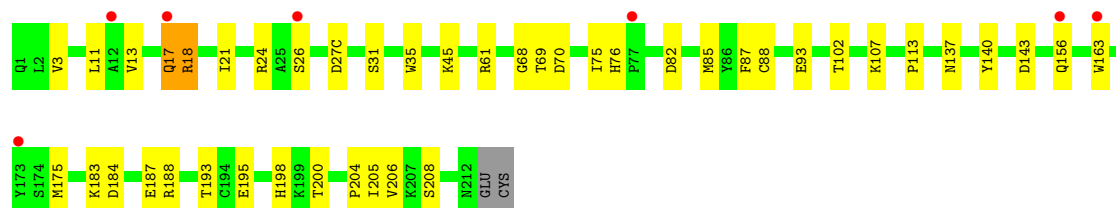
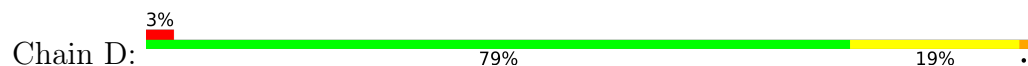




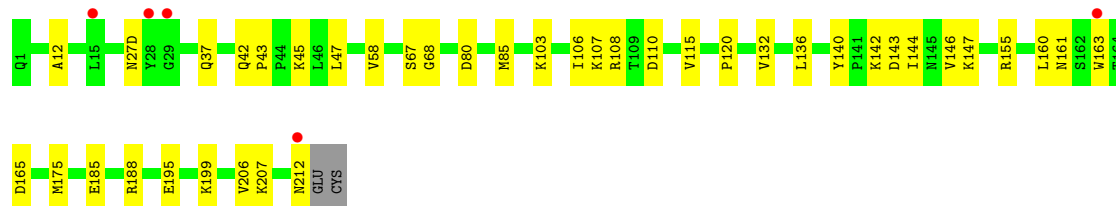
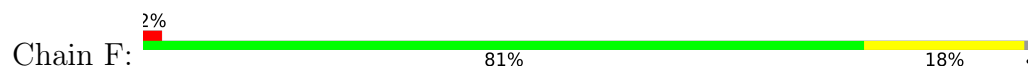
• Molecule 2: G2 Fab Light chain



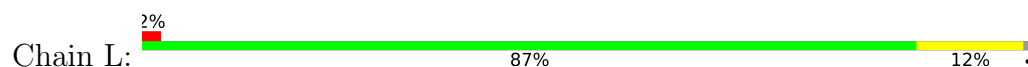
• Molecule 2: G2 Fab Light chain



• Molecule 2: G2 Fab Light chain



• Molecule 2: G2 Fab Light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.64Å 207.14Å 103.36Å 90.00° 97.70° 90.00°	Depositor
Resolution (Å)	38.41 – 2.10 38.41 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.2 (38.41-2.10) 97.3 (38.41-2.10)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.181 , 0.225 0.184 , 0.225	Depositor DCC
R_{free} test set	4773 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.197	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.1	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	0.96	EDS
Total number of atoms	13415	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 70.53 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.0748e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.45	0/1612	0.60	0/2203
1	C	0.40	0/1621	0.59	0/2215
1	E	0.38	0/1612	0.61	0/2203
1	H	0.40	0/1612	0.58	0/2203
2	B	0.45	0/1696	0.63	0/2304
2	D	0.38	0/1696	0.62	1/2304 (0.0%)
2	F	0.35	0/1696	0.56	0/2304
2	L	0.37	0/1696	0.55	0/2304
All	All	0.40	0/13241	0.59	1/18040 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	137	ASN	N-CA-C	6.15	119.72	109.95

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	1570	0	1524	24	0
1	C	1579	0	1535	15	0
1	E	1570	0	1524	25	1
1	H	1570	0	1524	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1659	0	1587	23	2
2	D	1659	0	1589	32	1
2	F	1659	0	1587	32	2
2	L	1659	0	1587	30	0
3	A	92	0	0	6	0
3	B	80	0	0	5	0
3	C	89	0	0	4	0
3	D	48	0	0	1	0
3	E	25	0	0	1	0
3	F	42	0	0	11	0
3	H	65	0	0	5	0
3	L	49	0	0	3	1
All	All	13415	0	12457	192	4

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 (192) 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:L:163:TRP:HE1	2:L:175:MET:CE	1.43	1.30
2:L:163:TRP:HE1	2:L:175:MET:HE2	1.11	1.11
2:L:163:TRP:NE1	2:L:175:MET:HE2	1.64	1.11
2:L:163:TRP:NE1	2:L:175:MET:CE	2.15	1.09
2:D:61:ARG:NH1	2:D:82:ASP:OD2	1.95	0.98
2:L:163:TRP:HE1	2:L:175:MET:HE3	1.27	0.96
1:C:41:PRO:O	3:C:301:HOH:O	1.87	0.92
2:F:45:LYS:NZ	3:F:304:HOH:O	2.10	0.84
1:H:2:VAL:N	3:H:303:HOH:O	2.09	0.84
2:L:2:LEU:HD11	2:L:93:GLU:CD	2.04	0.83
2:B:112:ALA:HB2	2:B:200:THR:HG21	1.57	0.83
1:A:184:PRO:O	1:A:187:THR:HG22	1.82	0.79
2:L:163:TRP:NE1	2:L:175:MET:HE3	1.91	0.79
2:F:155:ARG:NH2	2:F:185:GLU:OE2	2.15	0.78
1:A:2:VAL:N	3:A:303:HOH:O	2.18	0.76
2:F:212:ASN:O	3:F:301:HOH:O	2.02	0.76
2:B:146:VAL:O	3:B:301:HOH:O	2.04	0.74
2:F:108:ARG:O	3:F:302:HOH:O	2.05	0.74
1:E:202:SER:O	1:H:208:LYS:NZ	2.22	0.72
2:F:110:ASP:OD2	2:F:199:LYS:NZ	2.23	0.72
1:H:171:GLN:OE1	3:H:301:HOH:O	2.08	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:167:PRO:O	3:H:302:HOH:O	2.09	0.71
2:L:2:LEU:CD1	2:L:93:GLU:CD	2.63	0.70
2:F:185:GLU:OE1	3:F:303:HOH:O	2.08	0.70
1:C:79:TYR:OH	3:C:302:HOH:O	2.09	0.69
2:L:110:ASP:OD1	3:L:301:HOH:O	2.10	0.69
1:C:19:LYS:NZ	3:C:305:HOH:O	2.26	0.68
2:L:33:MET:HE2	2:L:88:CYS:HB2	1.76	0.68
1:A:126:PRO:HA	1:A:213:ARG:HH21	1.59	0.67
2:B:112:ALA:CB	2:B:200:THR:HG21	2.24	0.67
1:A:112:SER:O	3:A:301:HOH:O	2.13	0.66
1:E:39:GLN:HB2	1:E:45:LEU:HD13	1.77	0.66
1:E:37:LEU:HD13	1:E:47:TRP:HA	1.79	0.65
2:F:43:PRO:HD3	3:F:314:HOH:O	1.96	0.65
2:B:212:ASN:O	3:B:302:HOH:O	2.14	0.65
1:C:72:ASP:OD1	1:C:74:SER:OG	2.08	0.65
2:D:18:ARG:NH2	2:D:76:HIS:CE1	2.65	0.64
1:E:51:ILE:HD13	1:E:71:VAL:HG13	1.79	0.64
1:E:204:THR:HG22	1:H:208:LYS:HD2	1.80	0.64
2:F:107:LYS:HD2	2:F:140:TYR:OH	1.97	0.64
1:C:51:ILE:HD13	1:C:71:VAL:HG23	1.78	0.64
2:L:2:LEU:HD11	2:L:93:GLU:OE1	1.97	0.64
2:F:42:GLN:HG2	2:F:43:PRO:HD2	1.79	0.63
2:F:195:GLU:HG3	2:F:206:VAL:HG12	1.80	0.63
1:A:82:PHE:HD2	1:A:82(C):LEU:HD21	1.63	0.63
1:H:11:LEU:HD22	1:H:147:PRO:HG3	1.82	0.62
1:A:7:SER:OG	3:A:302:HOH:O	2.16	0.62
2:F:165:ASP:HB3	3:F:311:HOH:O	1.99	0.62
1:H:51:ILE:HD13	1:H:71:VAL:HG13	1.82	0.62
2:L:103:LYS:NZ	3:L:302:HOH:O	2.28	0.62
1:E:37:LEU:CD1	1:E:47:TRP:HA	2.30	0.61
1:E:187:THR:O	1:E:187:THR:HG22	1.99	0.61
2:D:193:THR:HG23	2:D:208:SER:HB3	1.84	0.60
1:A:209:LYS:HB3	1:C:203:SER:HB3	1.84	0.60
1:H:208:LYS:HE2	2:L:123:GLU:OE2	2.02	0.60
1:H:4:LEU:HD22	1:H:24:THR:HG22	1.82	0.59
2:B:180:THR:HG23	3:B:308:HOH:O	2.03	0.59
2:B:155:ARG:HD2	3:B:306:HOH:O	2.02	0.59
2:F:107:LYS:O	3:F:305:HOH:O	2.17	0.58
2:B:123:GLU:HG2	1:C:118:PRO:HD3	1.86	0.58
2:D:61:ARG:HH12	2:D:82:ASP:CG	2.11	0.58
1:E:66:LYS:NZ	1:E:82(A):SER:O	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:2:VAL:N	3:E:304:HOH:O	2.36	0.57
2:L:163:TRP:CE2	2:L:175:MET:HE2	2.38	0.57
1:E:87:SER:OG	1:E:110:THR:HA	2.05	0.57
2:F:12:ALA:HB1	2:F:107:LYS:HD3	1.86	0.56
1:C:19:LYS:HD3	1:C:81:GLN:HG2	1.86	0.56
2:F:85:MET:HE3	2:F:103:LYS:HB2	1.88	0.56
2:F:12:ALA:CB	2:F:107:LYS:HD3	2.36	0.56
2:L:126:THR:HG22	3:L:321:HOH:O	2.05	0.55
2:D:18:ARG:HH21	2:D:76:HIS:CE1	2.24	0.55
1:E:12:VAL:O	1:E:111:VAL:HA	2.07	0.55
1:A:169:VAL:HG21	2:B:160:LEU:HD11	1.87	0.55
2:B:160:LEU:C	2:B:160:LEU:HD13	2.31	0.55
1:A:169:VAL:HG11	2:B:160:LEU:HD21	1.88	0.54
1:A:171:GLN:HG3	3:A:319:HOH:O	2.06	0.54
2:L:13:VAL:HG21	2:L:19:ALA:HB2	1.89	0.54
2:D:18:ARG:NH2	2:D:76:HIS:CG	2.76	0.53
1:A:100(A):MET:HE2	1:A:103:TRP:CZ2	2.44	0.53
2:L:163:TRP:CD1	2:L:175:MET:HE3	2.44	0.53
2:D:27(C):ASP:OD1	2:D:31:SER:OG	2.17	0.53
1:H:126:PRO:HA	1:H:213:ARG:HH21	1.73	0.53
2:L:163:TRP:CD1	2:L:175:MET:HG3	2.44	0.53
2:D:69:THR:OG1	2:D:70:ASP:OD1	2.24	0.52
2:D:11:LEU:HD11	2:D:13:VAL:HG23	1.91	0.52
1:H:13:LYS:H	1:H:13:LYS:HD2	1.74	0.52
2:L:183:LYS:O	2:L:187:GLU:HG3	2.09	0.51
1:E:183:VAL:HG22	1:E:184:PRO:HD2	1.92	0.51
1:A:61:GLN:NE2	3:A:309:HOH:O	2.34	0.51
2:D:45:LYS:HD2	3:D:313:HOH:O	2.11	0.50
2:B:2:LEU:HD21	2:B:27:GLU:HB3	1.94	0.50
2:B:145:ASN:HB2	2:B:197:THR:CG2	2.41	0.50
2:D:18:ARG:NH2	2:D:76:HIS:ND1	2.59	0.50
2:D:195:GLU:HG3	2:D:206:VAL:HG22	1.94	0.50
2:L:1:GLN:OE1	2:L:1:GLN:N	2.26	0.50
2:D:93:GLU:OE2	2:D:93:GLU:HA	2.12	0.49
2:F:43:PRO:CD	3:F:314:HOH:O	2.58	0.49
1:E:46:GLU:OE2	1:E:62:LYS:HE2	2.12	0.49
2:D:61:ARG:NH1	2:D:82:ASP:CG	2.68	0.49
2:F:161:ASN:HB3	2:F:175:MET:HE3	1.94	0.49
2:B:145:ASN:HB2	2:B:197:THR:HG23	1.94	0.49
2:L:27(C):ASP:OD1	2:L:31:SER:OG	2.24	0.49
1:A:37:LEU:HD21	1:A:100(A):MET:HE1	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:117:THR:OG1	2:L:126:THR:HG21	2.13	0.49
2:L:33:MET:HE3	2:L:89:GLN:C	2.38	0.49
2:F:47:LEU:HA	2:F:58:VAL:HG21	1.93	0.48
2:B:39:LYS:HE3	3:B:351:HOH:O	2.13	0.48
1:E:12:VAL:CG2	1:E:111:VAL:HG12	2.44	0.48
1:E:83:THR:HG22	1:E:85:GLU:H	1.78	0.48
2:F:136:LEU:HD21	2:F:146:VAL:HG12	1.95	0.48
1:H:19:LYS:HE2	1:H:79:TYR:CD1	2.49	0.48
1:C:57:THR:O	3:C:304:HOH:O	2.20	0.48
1:A:98:SER:HB3	2:B:32:PHE:CE1	2.49	0.48
1:C:11:LEU:HD23	1:C:110:THR:HB	1.96	0.48
2:D:18:ARG:HA	2:D:75:ILE:O	2.13	0.48
2:D:3:VAL:H	2:D:26:SER:HB3	1.78	0.47
2:D:163:TRP:CD1	2:D:175:MET:HG3	2.49	0.47
1:H:50:TRP:CD1	1:H:50:TRP:C	2.93	0.47
2:D:195:GLU:HG2	2:D:204:PRO:HB2	1.96	0.47
1:H:103:TRP:CE3	2:L:44:PRO:HD2	2.50	0.47
2:D:85:MET:HG2	2:D:87:PHE:CZ	2.49	0.47
2:F:115:VAL:O	2:F:207:LYS:HE2	2.15	0.47
2:F:120:PRO:HD3	2:F:132:VAL:HG22	1.97	0.47
1:A:4:LEU:O	3:A:304:HOH:O	2.21	0.46
2:F:185:GLU:HG3	2:F:188:ARG:NH1	2.30	0.46
1:A:37:LEU:HD11	1:A:100(A):MET:HE3	1.98	0.46
1:E:50:TRP:C	1:E:50:TRP:CD1	2.93	0.46
2:F:37:GLN:OE1	2:F:45:LYS:HE2	2.15	0.46
2:F:147:LYS:HB3	2:F:147:LYS:HE2	1.31	0.46
2:B:27(D):ASN:HB3	2:B:32:PHE:HE2	1.81	0.45
1:E:12:VAL:HG23	1:E:111:VAL:HG12	1.96	0.45
1:C:29:PHE:CD2	1:C:76:SER:HA	2.52	0.45
2:F:103:LYS:NZ	3:F:311:HOH:O	2.44	0.45
1:A:37:LEU:HD11	1:A:100(A):MET:CE	2.47	0.45
2:D:35:TRP:CZ3	2:D:88:CYS:HB3	2.52	0.45
2:F:142:LYS:HE3	2:F:143:ASP:OD1	2.17	0.45
2:F:160:LEU:HD23	2:F:160:LEU:HA	1.85	0.45
1:H:20:LEU:HD22	1:H:107:THR:HG21	1.99	0.45
1:A:93:ALA:HB3	1:A:100(A):MET:HE3	1.98	0.44
2:D:107:LYS:HA	2:D:140:TYR:OH	2.17	0.44
1:A:18:VAL:HG22	1:A:82(C):LEU:HD11	1.99	0.44
2:B:160:LEU:HD13	2:B:161:ASN:N	2.33	0.44
1:A:98:SER:HB3	2:B:32:PHE:HE1	1.83	0.44
1:A:103:TRP:CE3	2:B:44:PRO:HD2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:THR:HG23	1:A:209:LYS:HE2	2.00	0.44
2:F:43:PRO:N	3:F:314:HOH:O	2.51	0.44
3:H:301:HOH:O	2:L:160:LEU:HD23	2.17	0.44
2:L:1:GLN:H3	2:L:1:GLN:CD	2.16	0.44
2:D:183:LYS:O	2:D:187:GLU:HG3	2.17	0.44
1:E:38:LYS:HB2	1:E:48:ILE:HD11	1.99	0.44
1:H:199:HIS:ND1	1:H:202:SER:OG	2.45	0.44
1:H:159:LEU:HD13	1:H:181:VAL:HG21	1.99	0.44
1:E:146:PHE:HB2	1:E:174:LEU:HD23	2.00	0.44
2:L:2:LEU:HD13	2:L:93:GLU:CD	2.42	0.44
2:D:21:ILE:HG12	2:D:102:THR:HG21	1.99	0.43
1:H:7:SER:OG	3:H:304:HOH:O	2.21	0.43
1:A:66:LYS:NZ	1:A:86:ASP:OD2	2.41	0.43
1:H:4:LEU:HD22	1:H:24:THR:CG2	2.47	0.43
1:C:187:THR:O	1:C:191:GLU:HB2	2.18	0.43
1:C:154:TRP:CZ3	1:C:195:CYS:HB3	2.53	0.43
2:B:33:MET:HG2	2:B:71:PHE:CG	2.54	0.43
1:E:119:PRO:HB3	1:E:145:TYR:HB3	2.00	0.43
2:D:198:HIS:HB3	2:D:200:THR:HG22	2.01	0.42
1:E:67:ALA:HA	1:E:81:GLN:O	2.19	0.42
1:C:19:LYS:HB2	1:C:81:GLN:HE21	1.85	0.42
2:F:144:ILE:HG22	2:F:163:TRP:CZ3	2.54	0.42
1:H:171:GLN:O	1:H:171:GLN:HG3	2.20	0.42
2:L:120:PRO:HG3	2:L:130:ALA:HB1	2.00	0.42
1:H:83:THR:OG1	1:H:84:THR:N	2.53	0.42
2:D:24:ARG:NH2	2:D:70:ASP:OD2	2.53	0.42
2:L:16:GLY:HA2	2:L:77:PRO:HB2	2.00	0.42
2:L:76:HIS:HA	2:L:77:PRO:HA	1.93	0.42
2:F:80:ASP:OD1	2:F:106:ILE:HG13	2.20	0.41
1:E:148:GLU:HG3	1:E:175:TYR:CE2	2.54	0.41
2:D:70:ASP:OD1	2:D:70:ASP:N	2.52	0.41
1:C:82:PHE:HD2	1:C:82(C):LEU:HD21	1.85	0.41
1:H:50:TRP:CH2	1:H:97:SER:HA	2.56	0.41
2:B:25:ALA:C	2:B:27:GLU:H	2.28	0.41
1:E:24:THR:HB	1:E:27:PHE:CE1	2.56	0.41
1:H:13:LYS:HD2	1:H:13:LYS:N	2.36	0.41
2:D:13:VAL:HG13	2:D:17:GLN:HG2	2.03	0.41
2:D:184:ASP:O	2:D:188:ARG:HG3	2.21	0.41
2:B:105:GLU:OE2	2:B:142:LYS:HE3	2.21	0.41
2:D:18:ARG:HH21	2:D:76:HIS:CG	2.39	0.41
2:D:113:PRO:HG2	2:D:205:ILE:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:85:MET:CE	2:F:103:LYS:HB2	2.51	0.41
2:F:43:PRO:HG2	3:F:319:HOH:O	2.21	0.40
1:A:46:GLU:OE2	1:A:62:LYS:HE3	2.22	0.40
2:B:1:GLN:OE1	2:B:1:GLN:N	2.45	0.40
2:D:24:ARG:CZ	2:D:70:ASP:OD2	2.69	0.40
1:E:72:ASP:CG	1:E:75:SER:HB3	2.47	0.40
2:D:18:ARG:NH2	2:D:76:HIS:CD2	2.90	0.40

All (4) 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)
2:B:107:LYS:NZ	2:F:67:SER:OG[2_354]	1.99	0.21
2:D:156:GLN:NE2	1:E:75:SER:OG[1_556]	2.04	0.16
3:L:324:HOH:O	3:L:335:HOH:O[1_455]	2.09	0.11
2:B:18:ARG:NH1	2:F:27(D):ASN:OD1[2_354]	2.10	0.10

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	206/229 (90%)	201 (98%)	5 (2%)	0	100	100
1	C	207/229 (90%)	201 (97%)	6 (3%)	0	100	100
1	E	206/229 (90%)	195 (95%)	11 (5%)	0	100	100
1	H	206/229 (90%)	200 (97%)	6 (3%)	0	100	100
2	B	214/218 (98%)	209 (98%)	4 (2%)	1 (0%)	24	22
2	D	214/218 (98%)	209 (98%)	4 (2%)	1 (0%)	24	22
2	F	214/218 (98%)	208 (97%)	5 (2%)	1 (0%)	24	22
2	L	214/218 (98%)	209 (98%)	4 (2%)	1 (0%)	24	22

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1681/1788 (94%)	1632 (97%)	45 (3%)	4 (0%)	43	44

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	68	GLY
2	L	68	GLY
2	D	68	GLY
2	F	68	GLY

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	180/194 (93%)	180 (100%)	0	100	100
1	C	181/194 (93%)	181 (100%)	0	100	100
1	E	180/194 (93%)	179 (99%)	1 (1%)	78	86
1	H	180/194 (93%)	180 (100%)	0	100	100
2	B	190/192 (99%)	190 (100%)	0	100	100
2	D	190/192 (99%)	187 (98%)	3 (2%)	55	64
2	F	190/192 (99%)	190 (100%)	0	100	100
2	L	190/192 (99%)	190 (100%)	0	100	100
All	All	1481/1544 (96%)	1477 (100%)	4 (0%)	86	91

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	17	GLN
2	D	18	ARG
2	D	143	ASP
1	E	192	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
2	B	38	GLN
2	B	157	ASN
2	B	190	ASN
1	C	3	GLN
1	C	39	GLN
1	C	81	GLN
1	C	164	HIS
2	D	17	GLN
2	D	38	GLN
2	D	49	HIS
1	E	39	GLN
2	F	38	GLN
2	F	49	HIS
2	F	74	ASN
2	F	76	HIS
2	L	53	ASN
2	L	54	GLN
2	L	157	ASN
2	L	190	ASN
2	L	210	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	210/229 (91%)	-0.04	3 (1%) 73 75	25, 47, 72, 83	0
1	C	211/229 (92%)	0.07	1 (0%) 87 88	27, 50, 80, 101	0
1	E	210/229 (91%)	0.64	14 (6%) 24 25	35, 67, 95, 108	0
1	H	210/229 (91%)	0.06	3 (1%) 73 75	31, 47, 70, 91	0
2	B	216/218 (99%)	-0.05	4 (1%) 66 69	27, 48, 82, 109	0
2	D	216/218 (99%)	0.44	7 (3%) 50 53	30, 61, 98, 120	0
2	F	216/218 (99%)	0.40	5 (2%) 61 64	39, 61, 91, 112	0
2	L	216/218 (99%)	0.18	5 (2%) 61 64	33, 56, 80, 109	0
All	All	1705/1788 (95%)	0.21	42 (2%) 58 61	25, 54, 88, 120	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	29	GLY	3.4
1	A	64	THR	3.3
1	E	5	GLN	3.3
2	F	163	TRP	3.2
2	D	173	TYR	3.2
2	D	163	TRP	3.0
1	E	187	THR	3.0
1	E	107	THR	2.8
1	A	2	VAL	2.8
1	E	108	SER	2.7
1	H	64	THR	2.6
2	L	2	LEU	2.6
1	E	185	SER	2.5
1	H	4	LEU	2.5
1	E	191	GLU	2.5
2	D	17	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
2	D	26	SER	2.5
1	A	96	GLY	2.5
1	E	85	GLU	2.4
1	E	170	LEU	2.4
1	C	71	VAL	2.4
2	D	77	PRO	2.4
2	F	28	TYR	2.3
2	D	12	ALA	2.3
1	E	9	GLY	2.3
1	E	172	SER	2.3
2	B	27	GLU	2.3
2	B	81	ASP	2.3
2	F	212	ASN	2.2
1	E	11	LEU	2.2
2	B	28	TYR	2.2
1	H	2	VAL	2.1
2	L	67	SER	2.1
2	L	160	LEU	2.1
2	L	81	ASP	2.1
1	E	111	VAL	2.1
2	B	160	LEU	2.1
2	D	156	GLN	2.0
2	L	126	THR	2.0
1	E	65	GLY	2.0
1	E	174	LEU	2.0
2	F	15	LEU	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.