



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:41 PM UTC

PDB ID : 5CP7 / pdb_00005cp7
Title : Crystal Structure of an Antigen-Binding Fragment of Monoclonal Antibody against Sulfonamides
Authors : Wang, Z.; Shen, J.; Li, C.; Li, Y.; Wen, K.; Yu, X.; Zhang, X.
Deposited on : 2015-07-21
Resolution : 3.01 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

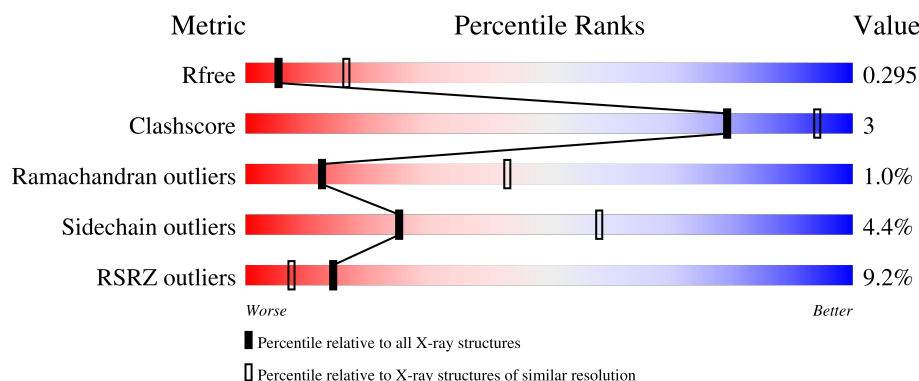
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 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	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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	92% 7% .
1	C	218	87% 12% .
1	E	218	17% 90% 8% .
1	G	218	6% 94% 6%
2	B	214	3% 92% 7% .

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Mol	Chain	Length	Quality of chain
2	D	214	13% 85% 10% 5%
2	F	214	32% 87% 9%
2	H	214	2% 83% 15%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13178 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 Light Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	0	1	0
			1700	1061	288	345	6			
1	C	218	Total	C	N	O	S	0	1	0
			1702	1062	288	346	6			
1	E	218	Total	C	N	O	S	0	0	0
			1686	1052	285	343	6			
1	G	218	Total	C	N	O	S	0	1	0
			1702	1062	289	345	6			

- Molecule 2 is a protein called Heavy Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	214	Total	C	N	O	S	0	0	0
			1628	1027	272	323	6			
2	D	204	Total	C	N	O	S	0	1	0
			1573	999	260	308	6			
2	F	207	Total	C	N	O	S	0	0	0
			1576	995	264	311	6			
2	H	212	Total	C	N	O	S	0	0	0
			1610	1017	268	319	6			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	O	0	0
			1	1		

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: Light Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7

Chain A: 



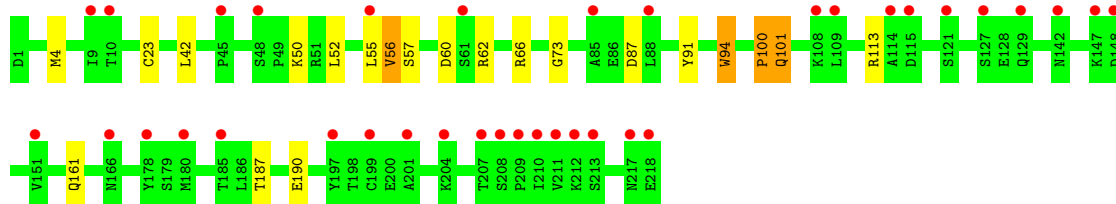
- Molecule 1: Light Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7

Chain C: 

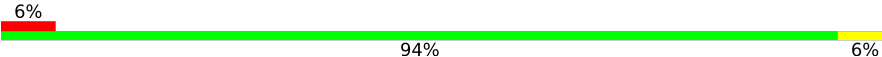


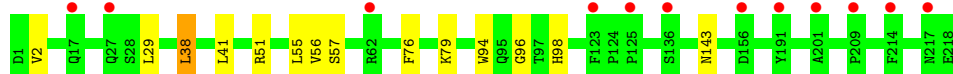
- Molecule 1: Light Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7

Chain E: 

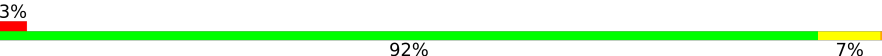


- Molecule 1: Light Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7

Chain G: 

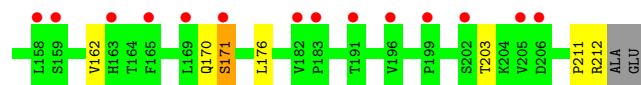
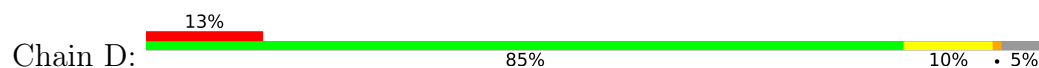


- Molecule 2: Heavy Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7

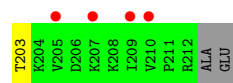
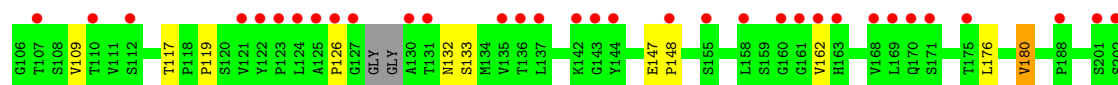
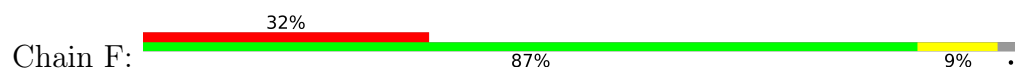
Chain B: 



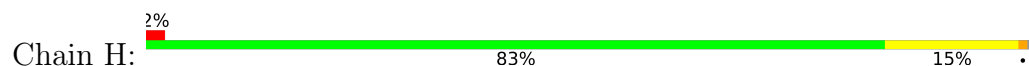
- Molecule 2: Heavy Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7



- Molecule 2: Heavy Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7



- Molecule 2: Heavy Chain of Antigen-Binding Fragment of Monoclonal Antibody of 4C7



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	176.10Å 96.51Å 138.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.88 – 3.01 29.88 – 3.01	Depositor EDS
% Data completeness (in resolution range)	94.6 (29.88-3.01) 94.6 (29.88-3.01)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.42 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.260 , 0.302 0.257 , 0.295	Depositor DCC
R_{free} test set	2275 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	41.6	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 20.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	13178	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.55% 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	0.42	0/1735	0.66	0/2356
1	C	0.41	0/1737	0.64	0/2357
1	E	0.45	0/1721	0.66	0/2338
1	G	0.43	0/1737	0.63	0/2357
2	B	0.42	0/1675	0.68	0/2297
2	D	0.42	0/1619	0.68	2/2220 (0.1%)
2	F	0.47	0/1622	0.70	1/2224 (0.0%)
2	H	0.43	0/1655	0.66	0/2270
All	All	0.43	0/13501	0.67	3/18419 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	32	GLY	N-CA-C	7.91	117.41	110.21
2	D	13	THR	CA-C-N	5.05	126.16	119.84
2	D	13	THR	C-N-CA	5.05	126.16	119.84

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	1700	0	1655	8	0
1	C	1702	0	1659	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1686	0	1634	11	0
1	G	1702	0	1661	6	0
2	B	1628	0	1566	10	0
2	D	1573	0	1515	10	0
2	F	1576	0	1506	5	0
2	H	1610	0	1549	12	0
3	D	1	0	0	0	0
All	All	13178	0	12745	69	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 3.

All (69) 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:A:55:LEU:O	1:A:57:SER:N	2.23	0.70
1:C:55:LEU:O	1:C:57:SER:N	2.25	0.69
1:G:55:LEU:O	1:G:57:SER:N	2.28	0.66
2:H:14:PRO:O	2:H:86:VAL:O	2.14	0.66
1:E:55:LEU:O	1:E:57:SER:N	2.30	0.64
1:C:91:TYR:CE1	1:C:109:LEU:HD12	2.33	0.64
2:F:18:LEU:HD11	2:F:109:VAL:HG11	1.85	0.58
2:H:3:GLN:C	2:H:4:LEU:HD23	2.29	0.57
1:G:41:LEU:HD23	1:G:51:ARG:HA	1.88	0.54
2:H:91:THR:HG23	2:H:110:THR:HA	1.88	0.54
2:B:41:PHE:O	2:B:43:GLY:O	2.26	0.53
2:H:82:GLN:NE2	2:H:84:ASN:OD1	2.41	0.52
2:D:30:THR:HG21	2:D:74:THR:HG23	1.92	0.52
2:B:91:THR:HG23	2:B:110:THR:HA	1.93	0.51
2:F:147:GLU:HB3	2:F:148:PRO:HA	1.93	0.50
1:E:94:TRP:CD1	1:E:94:TRP:C	2.89	0.50
2:B:3:GLN:C	2:B:4:LEU:HD12	2.37	0.50
2:B:176:LEU:C	2:B:176:LEU:HD12	2.37	0.49
2:D:4:LEU:HD22	2:D:24:VAL:HG12	1.94	0.49
1:E:66:ARG:NH2	1:E:87:ASP:OD2	2.43	0.49
1:E:4:MET:HE3	1:E:23:CYS:SG	2.54	0.48
2:D:91:THR:HG23	2:D:110:THR:HA	1.95	0.48
2:F:162:VAL:O	2:F:162:VAL:HG13	2.13	0.48
1:C:42:LEU:HD13	1:C:91:TYR:CE1	2.49	0.47
1:A:84:GLU:HG2	2:B:31:SER:CB	2.46	0.46
2:D:32:GLY:O	2:D:54:TYR:CD1	2.69	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:38:LEU:HD22	1:G:76:PHE:CG	2.51	0.46
1:C:38:LEU:HD13	1:C:76:PHE:CD1	2.51	0.46
1:E:42:LEU:HB2	1:E:52:LEU:HD11	1.97	0.46
1:G:29:LEU:HD12	1:G:76:PHE:CE1	2.51	0.45
1:C:83:VAL:HG23	1:C:83:VAL:O	2.15	0.45
2:D:176:LEU:C	2:D:176:LEU:HD12	2.41	0.45
1:A:84:GLU:HG2	2:B:31:SER:HB3	1.98	0.45
1:C:166:ASN:HB3	1:C:180:MET:HE2	1.99	0.45
1:E:55:LEU:O	1:E:56:VAL:C	2.59	0.45
2:B:64:LEU:O	2:B:65:ARG:C	2.60	0.45
2:D:49:MET:HE1	2:D:81:LEU:HD21	1.99	0.45
1:E:187:THR:HB	1:E:190:GLU:HB2	1.99	0.45
2:F:176:LEU:C	2:F:176:LEU:HD12	2.43	0.44
2:H:151:VAL:HG11	2:H:178:SER:HB3	2.00	0.44
2:B:42:PRO:C	2:B:43:GLY:O	2.58	0.43
2:H:4:LEU:HD22	2:H:24:VAL:HB	2.00	0.43
1:E:50:LYS:NZ	1:E:60:ASP:OD2	2.51	0.43
1:C:194:HIS:O	1:C:216:ARG:NH1	2.52	0.42
1:A:38:LEU:HD22	1:A:76:PHE:CG	2.54	0.42
2:B:14:PRO:O	2:B:86:VAL:O	2.37	0.42
2:H:43:GLY:O	2:H:44:ASN:HB2	2.20	0.42
2:B:52:ILE:HG23	2:B:52:ILE:O	2.19	0.42
1:G:2:VAL:HG21	1:G:98:HIS:HB2	2.02	0.42
2:D:170:GLN:O	2:D:171:SER:C	2.63	0.42
2:F:119:PRO:HD2	2:F:203:THR:HG21	2.00	0.42
2:D:3:GLN:C	2:D:4:LEU:HD23	2.44	0.42
2:H:151:VAL:HG11	2:H:178:SER:CB	2.49	0.42
1:E:42:LEU:HD13	1:E:91:TYR:CZ	2.55	0.41
2:H:137:LEU:HD12	2:H:192:VAL:HG11	2.02	0.41
1:A:155:ILE:HD11	1:A:184:LEU:HD21	2.01	0.41
1:A:38:LEU:HD22	1:A:76:PHE:CD2	2.54	0.41
1:E:62:ARG:O	1:E:62:ARG:HG3	2.19	0.41
1:A:110:GLU:HG3	1:A:178:TYR:OH	2.21	0.41
1:G:94:TRP:CH2	1:G:96:GLY:HA3	2.55	0.41
2:H:119:PRO:HB3	2:H:144:TYR:HB3	2.03	0.41
1:E:100:PRO:O	1:E:101:GLN:C	2.64	0.41
1:C:56:VAL:HG23	1:C:57:SER:N	2.36	0.40
2:D:211:PRO:O	2:D:212:ARG:HB3	2.21	0.40
1:A:29:LEU:HD12	1:A:76:PHE:CE1	2.56	0.40
1:C:13:VAL:HG21	1:C:83:VAL:HG11	2.01	0.40
2:D:4:LEU:HD12	2:D:97:ALA:HA	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:4:LEU:HD12	2:H:97:ALA:HA	2.03	0.40
2:H:176:LEU:HD12	2:H:176:LEU:C	2.46	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	217/218 (100%)	209 (96%)	7 (3%)	1 (0%)	24	59
1	C	217/218 (100%)	208 (96%)	8 (4%)	1 (0%)	24	59
1	E	216/218 (99%)	197 (91%)	15 (7%)	4 (2%)	6	28
1	G	217/218 (100%)	204 (94%)	11 (5%)	2 (1%)	14	46
2	B	212/214 (99%)	195 (92%)	16 (8%)	1 (0%)	24	59
2	D	201/214 (94%)	184 (92%)	15 (8%)	2 (1%)	12	43
2	F	203/214 (95%)	178 (88%)	21 (10%)	4 (2%)	6	26
2	H	210/214 (98%)	189 (90%)	19 (9%)	2 (1%)	12	43
All	All	1693/1728 (98%)	1564 (92%)	112 (7%)	17 (1%)	12	43

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	VAL
1	C	56	VAL
1	E	56	VAL
1	E	101	GLN
1	G	56	VAL
2	B	65	ARG
2	D	14	PRO
2	D	171	SER

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Mol	Chain	Res	Type
2	F	132	ASN
2	F	133	SER
2	H	160	GLY
2	F	180	VAL
1	G	143	ASN
2	H	131	THR
2	F	126	PRO
1	E	73	GLY
1	E	100	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	195/195 (100%)	186 (95%)	9 (5%)	24	57
1	C	196/195 (100%)	180 (92%)	16 (8%)	10	36
1	E	193/195 (99%)	190 (98%)	3 (2%)	55	78
1	G	196/195 (100%)	194 (99%)	2 (1%)	68	83
2	B	186/186 (100%)	180 (97%)	6 (3%)	34	66
2	D	181/186 (97%)	176 (97%)	5 (3%)	38	69
2	F	180/186 (97%)	171 (95%)	9 (5%)	22	54
2	H	185/186 (100%)	168 (91%)	17 (9%)	8	31
All	All	1512/1524 (99%)	1445 (96%)	67 (4%)	25	58

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	THR
1	A	38	LEU
1	A	62	ARG
1	A	86	GLU
1	A	109	LEU
1	A	110	GLU

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Mol	Chain	Res	Type
1	A	113	ARG
1	A	180	MET
1	A	211	VAL
2	B	2	VAL
2	B	31	SER
2	B	65	ARG
2	B	88	THR
2	B	110	THR
2	B	207	LYS
1	C	24	LYS
1	C	61	SER
1	C	62	ARG
1	C	68	THR
1	C	84	GLU
1	C	86	GLU
1	C	112	LYS
1	C	162	ASN
1	C	180	MET
1	C	181	SER
1	C	190[A]	GLU
1	C	190[B]	GLU
1	C	193	ARG
1	C	196	THR
1	C	212	LYS
1	C	217	ASN
2	D	6	GLU
2	D	24	VAL
2	D	68	ILE
2	D	162	VAL
2	D	203	THR
1	E	94	TRP
1	E	113	ARG
1	E	161	GLN
2	F	12	VAL
2	F	52	ILE
2	F	59	ASN
2	F	65	ARG
2	F	84	ASN
2	F	99	TYR
2	F	105	GLN
2	F	117	THR
2	F	180	VAL

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Mol	Chain	Res	Type
1	G	38	LEU
1	G	79	LYS
2	H	3	GLN
2	H	4	LEU
2	H	6	GLU
2	H	7	SER
2	H	12	VAL
2	H	24	VAL
2	H	45	LYS
2	H	65	ARG
2	H	76	LYS
2	H	89	GLU
2	H	139	CYS
2	H	168	VAL
2	H	184	SER
2	H	194	CYS
2	H	195	ASN
2	H	203	THR
2	H	207	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
2	B	40	GLN
2	B	78	GLN
2	B	84	ASN
2	B	105	GLN
2	B	198	HIS
1	C	17	GLN
1	C	43	GLN
2	D	16	GLN
2	D	36	HIS
1	E	47	GLN
1	E	101	GLN
1	E	142	ASN
1	E	161	GLN
1	E	166	ASN
1	E	194	HIS
1	E	195	ASN
2	F	40	GLN
2	F	59	ASN

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Mol	Chain	Res	Type
2	F	82	GLN
1	G	43	GLN
2	H	16	GLN
2	H	40	GLN
2	H	78	GLN
2	H	82	GLN
2	H	84	ASN
2	H	105	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	A	218/218 (100%)	0.07	1 (0%) 87 72	7, 16, 22, 31	1 (0%)
1	C	218/218 (100%)	0.13	1 (0%) 87 72	8, 18, 23, 26	1 (0%)
1	E	218/218 (100%)	1.30	36 (16%) 4 3	16, 44, 68, 72	0
1	G	218/218 (100%)	0.95	12 (5%) 30 15	13, 46, 59, 69	1 (0%)
2	B	214/214 (100%)	0.40	6 (2%) 55 32	14, 23, 41, 58	0
2	D	204/214 (95%)	1.02	28 (13%) 6 4	9, 32, 71, 84	1 (0%)
2	F	207/214 (96%)	1.63	69 (33%) 1 1	31, 52, 77, 80	0
2	H	212/214 (99%)	0.27	4 (1%) 66 43	6, 16, 36, 55	0
All	All	1709/1728 (98%)	0.72	157 (9%) 14 7	6, 26, 67, 84	4 (0%)

All (157) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	131	THR	5.9
1	E	199	CYS	4.9
2	F	127	GLY	4.7
2	H	1	ASP	4.2
2	D	202	SER	4.0
2	F	34	SER	4.0
2	F	161	GLY	3.8
1	E	85	ALA	3.7
2	H	130	ALA	3.6
2	F	142	LYS	3.5
2	F	130	ALA	3.5
2	F	4	LEU	3.5
1	E	109	LEU	3.5
2	F	32	GLY	3.5
2	F	163	HIS	3.4
2	B	128	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
2	D	163	HIS	3.3
2	D	171	SER	3.2
2	F	169	LEU	3.1
2	F	135	VAL	3.1
2	D	199	PRO	3.1
1	G	27[A]	GLN	3.1
2	D	105[A]	GLN	3.1
2	F	121	VAL	3.0
2	F	162	VAL	3.0
2	D	191	THR	3.0
2	F	124	LEU	2.9
2	F	16	GLN	2.9
1	E	207	THR	2.9
2	F	205	VAL	2.9
2	F	136	THR	2.8
2	F	168	VAL	2.8
2	B	32	GLY	2.8
2	F	160	GLY	2.8
2	F	170	GLN	2.8
2	D	159	SER	2.8
2	D	169	LEU	2.8
2	F	25	THR	2.8
2	D	134	MET	2.7
1	E	211	VAL	2.7
1	C	190[A]	GLU	2.7
2	D	206	ASP	2.7
2	D	158	LEU	2.7
1	E	180	MET	2.7
2	F	53	HIS	2.7
2	D	182	VAL	2.7
1	E	9	ILE	2.7
2	B	30	THR	2.6
2	F	102	TYR	2.6
2	D	183	PRO	2.6
2	H	132	ASN	2.6
1	E	147	LYS	2.6
1	E	55	LEU	2.6
1	E	201	ALA	2.6
2	F	64	LEU	2.5
2	D	11	LEU	2.5
2	F	137	LEU	2.5
1	E	209	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
2	F	33	TYR	2.5
2	F	123	PRO	2.5
2	D	68	ILE	2.5
2	F	24	VAL	2.5
2	F	87	THR	2.5
2	F	143	GLY	2.5
2	D	34	SER	2.4
2	F	15	SER	2.4
2	F	112	SER	2.4
1	E	166	ASN	2.4
2	B	131	THR	2.4
2	F	71	THR	2.4
1	E	197	TYR	2.4
1	E	127	SER	2.4
2	F	44	ASN	2.4
1	E	115	ASP	2.4
1	G	17	GLN	2.4
2	F	7	SER	2.4
1	G	125	PRO	2.4
1	E	212	LYS	2.4
2	D	137	LEU	2.4
2	F	158	LEU	2.4
2	F	19	SER	2.4
2	F	171	SER	2.4
2	D	116	THR	2.4
2	D	123	PRO	2.4
2	D	135	VAL	2.4
2	F	30	THR	2.4
1	E	204	LYS	2.4
2	D	65	ARG	2.4
2	F	68	ILE	2.3
1	E	210	ILE	2.3
2	B	130	ALA	2.3
1	E	178	TYR	2.3
1	G	217	ASN	2.3
1	E	151	VAL	2.3
2	D	196	VAL	2.3
1	G	209	PRO	2.3
2	F	101	GLY	2.3
2	F	209	ILE	2.3
1	E	148	ASP	2.3
2	F	52	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	181	SER	2.3
1	E	61	SER	2.3
2	F	202	SER	2.3
2	B	99	TYR	2.3
2	F	144	TYR	2.3
1	E	142	ASN	2.3
2	D	148	PRO	2.3
2	D	100	GLY	2.3
1	E	10	THR	2.3
2	F	107	THR	2.3
2	F	78	GLN	2.3
1	E	218	GLU	2.2
1	E	114	ALA	2.2
2	F	21	THR	2.2
1	E	108	LYS	2.2
1	E	121	SER	2.2
2	F	75	SER	2.2
2	D	126	PRO	2.2
2	F	125	ALA	2.2
2	F	126	PRO	2.2
2	F	54	TYR	2.2
1	E	48	SER	2.2
2	F	59	ASN	2.2
1	E	45	PRO	2.2
2	F	65	ARG	2.2
2	F	210	VAL	2.2
2	F	207	LYS	2.2
2	D	165	PHE	2.2
1	E	208	SER	2.2
1	G	136	SER	2.2
1	E	185	THR	2.1
2	F	122	TYR	2.1
2	F	31	SER	2.1
2	F	70	ILE	2.1
2	F	6	GLU	2.1
1	E	217	ASN	2.1
1	G	201	ALA	2.1
2	F	14	PRO	2.1
2	F	155	SER	2.1
1	G	156	ASP	2.1
2	F	175	THR	2.1
2	H	131	THR	2.1

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Mol	Chain	Res	Type	RSRZ
2	F	148	PRO	2.1
2	D	205	VAL	2.1
1	G	62	ARG	2.1
2	F	110	THR	2.1
2	D	32	GLY	2.1
1	E	213	SER	2.0
2	F	23	THR	2.0
1	E	129	GLN	2.0
1	G	214	PHE	2.0
2	F	201	SER	2.0
1	G	191	TYR	2.0
1	G	123	PHE	2.0
2	F	76	LYS	2.0
2	F	188	PRO	2.0
1	E	88	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.