



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:16 AM UTC

PDB ID : 4G8E / pdb_00004g8e
Title : Crystal Structure of clone18 TCR
Authors : Gras, S.; Bhati, M.; Rossjohn, J.
Deposited on : 2012-07-23
Resolution : 2.20 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

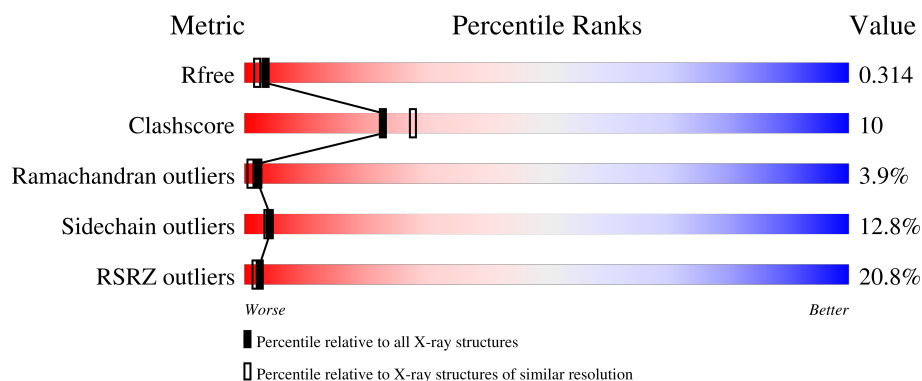
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	201	30% 64% 27% 7%
2	B	247	13% 67% 25% • • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3625 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 alpha chain clone 18 TCR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	201	Total	C	N	O	S	0	0	0
			1566	987	256	314	9			

- Molecule 2 is a protein called beta chain clone 18 TCR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	241	Total	C	N	O	S	2	0	0
			1913	1210	331	363	9			

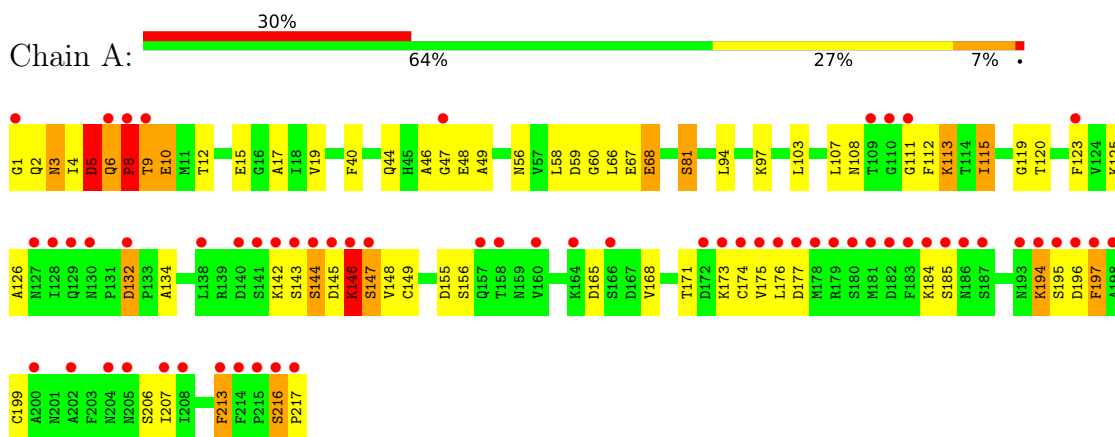
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	76	Total	O	0	0
			76	76		
3	B	70	Total	O	0	0
			70	70		

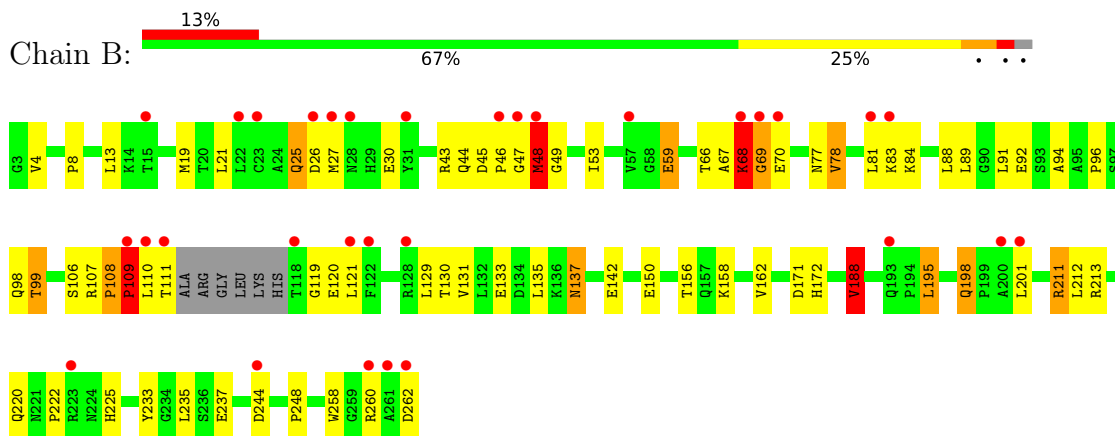
3 Residue-property plots

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: alpha chain clone 18 TCR



- Molecule 2: beta chain clone 18 TCR



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.42Å 84.09Å 58.28Å 90.00° 111.88° 90.00°	Depositor
Resolution (Å)	45.48 – 2.20 45.48 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.6 (45.48-2.20) 97.6 (45.48-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.19 (at 2.20Å)	Xtriage
Refinement program	BUSTER-TNT, BUSTER 2.10.0	Depositor
R, R_{free}	0.241 , 0.293 0.250 , 0.314	Depositor DCC
R_{free} test set	1062 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.465	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3625	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.86	0/1600	1.54	25/2165 (1.2%)
2	B	0.83	0/1964	1.35	13/2669 (0.5%)
All	All	0.84	0/3564	1.44	38/4834 (0.8%)

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	5	ASP	CA-CB-CG	11.65	124.25	112.60
1	A	17	ALA	N-CA-C	11.20	124.19	110.41
1	A	8	PRO	N-CA-C	9.03	131.07	112.47
1	A	10	GLU	N-CA-C	8.63	122.53	109.41
2	B	67	ALA	CA-C-N	8.18	137.17	121.54
2	B	67	ALA	C-N-CA	8.18	137.17	121.54
1	A	68	GLU	N-CA-C	7.87	120.55	111.11
1	A	206	SER	CA-C-N	7.49	128.69	121.65
1	A	206	SER	C-N-CA	7.49	128.69	121.65
2	B	108	PRO	N-CA-C	7.27	119.57	110.70
2	B	137	ASN	CA-CB-CG	7.10	119.70	112.60
2	B	47	GLY	CA-C-N	6.94	134.79	121.54
2	B	47	GLY	C-N-CA	6.94	134.79	121.54
2	B	68	LYS	N-CA-C	6.56	124.78	110.80
2	B	119	GLY	CA-C-N	6.53	130.15	120.87
2	B	119	GLY	C-N-CA	6.53	130.15	120.87
1	A	113	LYS	N-CA-C	6.46	120.13	108.69
1	A	3	ASN	N-CA-C	6.17	118.82	108.02
1	A	59	ASP	CA-CB-CG	5.95	118.55	112.60
1	A	197	PHE	CA-CB-CG	5.94	119.74	113.80
1	A	67	GLU	CA-C-N	5.94	128.55	120.54
1	A	67	GLU	C-N-CA	5.94	128.55	120.54
2	B	109	PRO	N-CA-C	5.92	124.66	112.47
1	A	4	ILE	CA-C-N	5.85	132.29	122.87
1	A	4	ILE	C-N-CA	5.85	132.29	122.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	GLY	CA-C-N	5.61	132.25	121.54
1	A	1	GLY	C-N-CA	5.61	132.25	121.54
2	B	67	ALA	N-CA-C	5.55	117.42	109.14
1	A	10	GLU	CA-C-N	5.44	131.75	121.69
1	A	10	GLU	C-N-CA	5.44	131.75	121.69
2	B	142	GLU	N-CA-C	-5.29	101.08	109.76
1	A	132	ASP	CA-CB-CG	5.27	117.87	112.60
1	A	143	SER	CA-C-N	5.14	131.36	121.54
1	A	143	SER	C-N-CA	5.14	131.36	121.54
2	B	188	VAL	N-CA-CB	5.13	119.07	110.86
1	A	146	LYS	CA-C-N	5.10	131.28	121.54
1	A	146	LYS	C-N-CA	5.10	131.28	121.54
1	A	213	PHE	CA-CB-CG	5.01	118.81	113.80

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	1566	0	1493	34	0
2	B	1913	0	1835	41	0
3	A	76	0	0	0	0
3	B	70	0	0	0	0
All	All	3625	0	3328	71	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 (71) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:48:MET:HB2	2:B:49:GLY:CA	1.82	1.08
2:B:48:MET:CB	2:B:49:GLY:HA2	1.88	1.03
2:B:48:MET:HB2	2:B:49:GLY:HA2	1.05	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:96:PRO:O	2:B:99:THR:HG23	1.80	0.81
1:A:149:CYS:HG	1:A:199:CYS:HG	0.81	0.81
1:A:44:GLN:HE22	2:B:44:GLN:HE22	1.29	0.79
1:A:56:ASN:HD21	1:A:81:SER:HB3	1.49	0.76
1:A:8:PRO:HG3	1:A:119:GLY:O	1.90	0.71
1:A:40:PHE:HE2	1:A:107:LEU:HB2	1.58	0.69
1:A:10:GLU:HG2	1:A:123:PHE:CE2	2.26	0.68
2:B:8:PRO:HD2	2:B:21:LEU:HD23	1.75	0.68
1:A:8:PRO:HA	1:A:9:THR:C	2.17	0.67
2:B:198:GLN:HB3	2:B:201:LEU:HD13	1.79	0.64
2:B:109:PRO:HA	2:B:111:THR:H	1.63	0.63
1:A:15:GLU:O	1:A:94:LEU:O	2.18	0.61
2:B:46:PRO:C	2:B:48:MET:H	2.06	0.60
2:B:45:ASP:O	2:B:48:MET:HB3	2.01	0.60
2:B:99:THR:HB	2:B:130:THR:HA	1.83	0.60
1:A:10:GLU:HG2	1:A:123:PHE:HE2	1.65	0.60
2:B:13:LEU:HD11	2:B:19:MET:HG2	1.85	0.57
2:B:107:ARG:HE	2:B:109:PRO:HG2	1.70	0.57
2:B:188:VAL:HG12	2:B:212:LEU:HD13	1.87	0.56
2:B:108:PRO:HD2	2:B:109:PRO:HD3	1.86	0.56
1:A:111:GLY:O	1:A:113:LYS:N	2.38	0.56
2:B:99:THR:HG22	2:B:131:VAL:H	1.70	0.55
1:A:5:ASP:C	1:A:8:PRO:HD2	2.31	0.55
1:A:8:PRO:HD3	1:A:120:THR:HA	1.88	0.55
1:A:165:ASP:HB3	1:A:168:VAL:HG12	1.88	0.55
2:B:109:PRO:HA	2:B:111:THR:N	2.22	0.54
1:A:12:THR:HG22	1:A:123:PHE:HD2	1.72	0.53
1:A:176:LEU:HD22	2:B:213:ARG:HB2	1.91	0.53
2:B:94:ALA:HA	2:B:98:GLN:NE2	2.25	0.52
1:A:40:PHE:CE2	1:A:107:LEU:HB2	2.43	0.51
1:A:145:ASP:C	1:A:147:SER:H	2.19	0.50
1:A:126:ALA:HB2	1:A:175:VAL:HG11	1.93	0.50
1:A:5:ASP:HB3	1:A:8:PRO:HD2	1.93	0.50
2:B:25:GLN:HG2	2:B:27:MET:H	1.76	0.49
2:B:13:LEU:HD21	2:B:19:MET:HB3	1.95	0.49
1:A:216:SER:H	1:A:217:PRO:HD2	1.78	0.48
1:A:6:GLN:N	1:A:8:PRO:HD2	2.28	0.48
2:B:99:THR:CG2	2:B:131:VAL:H	2.27	0.48
2:B:94:ALA:HA	2:B:98:GLN:HE22	1.79	0.48
2:B:21:LEU:HD12	2:B:89:LEU:HD23	1.95	0.47
1:A:56:ASN:ND2	1:A:81:SER:HB3	2.25	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:225:HIS:HB2	2:B:258:TRP:CZ3	2.50	0.47
1:A:175:VAL:HA	1:A:185:SER:O	2.14	0.47
2:B:30:GLU:HG2	2:B:84:LYS:HE3	1.96	0.46
1:A:144:SER:HB3	1:A:146:LYS:HE2	1.97	0.46
2:B:220:GLN:O	2:B:222:PRO:HD3	2.15	0.46
1:A:176:LEU:HD21	2:B:211:ARG:HB2	1.98	0.45
2:B:69:GLY:HA3	2:B:70:GLU:HA	1.47	0.45
1:A:134:ALA:HB2	1:A:213:PHE:HB3	1.99	0.44
2:B:172:HIS:HB3	2:B:233:TYR:HB2	1.99	0.43
1:A:125:LYS:HD2	1:A:156:SER:HB3	2.00	0.43
1:A:46:ALA:HA	1:A:47:GLY:HA2	1.82	0.43
2:B:235:LEU:HD22	2:B:248:PRO:HD2	1.99	0.43
2:B:77:ASN:HD21	2:B:92:GLU:HG3	1.83	0.43
2:B:45:ASP:HB3	2:B:46:PRO:CD	2.49	0.43
2:B:48:MET:CG	2:B:49:GLY:HA2	2.46	0.42
1:A:12:THR:HG22	1:A:123:PHE:CD2	2.53	0.42
2:B:107:ARG:HG3	2:B:120:GLU:O	2.18	0.42
1:A:108:ASN:HD21	1:A:115:ILE:HD13	1.84	0.42
2:B:46:PRO:HA	2:B:48:MET:HE2	2.02	0.42
2:B:78:VAL:HA	2:B:88:LEU:O	2.21	0.41
1:A:195:SER:N	1:A:196:ASP:HA	2.36	0.41
2:B:43:ARG:HB3	2:B:53:ILE:HD11	2.03	0.41
1:A:48:GLU:HG3	1:A:49:ALA:H	1.84	0.41
1:A:177:ASP:HA	1:A:184:LYS:HA	2.03	0.41
2:B:220:GLN:HA	2:B:260:ARG:O	2.21	0.41
1:A:171:THR:HG22	2:B:195:LEU:HD13	2.02	0.40
2:B:156:THR:HB	2:B:158:LYS:HE2	2.04	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	199/201 (99%)	175 (88%)	13 (6%)	11 (6%)	1	0
2	B	237/247 (96%)	220 (93%)	11 (5%)	6 (2%)	4	2
All	All	436/448 (97%)	395 (91%)	24 (6%)	17 (4%)	2	1

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	8	PRO
1	A	112	PHE
1	A	147	SER
1	A	173	LYS
1	A	194	LYS
2	B	48	MET
2	B	68	LYS
1	A	142	LYS
1	A	146	LYS
1	A	216	SER
2	B	59	GLU
2	B	109	PRO
1	A	60	GLY
1	A	144	SER
2	B	171	ASP
2	B	69	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	175/175 (100%)	155 (89%)	20 (11%)	5	5
2	B	208/212 (98%)	179 (86%)	29 (14%)	3	3
All	All	383/387 (99%)	334 (87%)	49 (13%)	4	4

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

Mol	Chain	Res	Type
1	A	3	ASN
1	A	5	ASP
1	A	6	GLN
1	A	8	PRO
1	A	9	THR
1	A	19	VAL
1	A	58	LEU
1	A	66	LEU
1	A	68	GLU
1	A	81	SER
1	A	97	LYS
1	A	103	LEU
1	A	115	ILE
1	A	132	ASP
1	A	148	VAL
1	A	155	ASP
1	A	174	CYS
1	A	194	LYS
1	A	197	PHE
1	A	207	ILE
2	B	4	VAL
2	B	25	GLN
2	B	26	ASP
2	B	48	MET
2	B	59	GLU
2	B	66	THR
2	B	68	LYS
2	B	78	VAL
2	B	81	LEU
2	B	83	LYS
2	B	91	LEU
2	B	99	THR
2	B	106	SER
2	B	109	PRO
2	B	110	LEU
2	B	121	LEU
2	B	129	LEU
2	B	133	GLU
2	B	135	LEU
2	B	137	ASN
2	B	150	GLU
2	B	162	VAL
2	B	188	VAL

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Mol	Chain	Res	Type
2	B	195	LEU
2	B	198	GLN
2	B	211	ARG
2	B	237	GLU
2	B	244	ASP
2	B	262	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	44	GLN
1	A	45	HIS
1	A	56	ASN
1	A	108	ASN
1	A	130	ASN
2	B	77	ASN
2	B	98	GLN
2	B	157	GLN
2	B	231	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	201/201 (100%)	1.50	61 (30%) 1 1	14, 37, 68, 83	12 (5%)
2	B	241/247 (97%)	0.85	31 (12%) 7 5	14, 31, 56, 81	2 (0%)
All	All	442/448 (98%)	1.15	92 (20%) 2 2	14, 33, 63, 83	14 (3%)

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	175	VAL	6.9
1	A	183	PHE	6.9
1	A	176	LEU	6.3
1	A	180	SER	6.2
1	A	8	PRO	5.8
1	A	184	LYS	5.4
1	A	195	SER	5.3
1	A	181	MET	5.1
1	A	174	CYS	5.0
1	A	217	PRO	5.0
2	B	118	THR	4.8
1	A	173	LYS	4.7
2	B	111	THR	4.3
2	B	110	LEU	4.3
1	A	111	GLY	4.3
2	B	69	GLY	4.3
1	A	182	ASP	4.2
1	A	123	PHE	4.0
1	A	9	THR	3.8
2	B	48	MET	3.7
2	B	109	PRO	3.6
2	B	70	GLU	3.5
1	A	130	ASN	3.4
1	A	213	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
2	B	46	PRO	3.2
1	A	207	ILE	3.2
1	A	145	ASP	3.2
1	A	109	THR	3.1
1	A	142	LYS	3.1
1	A	110	GLY	3.1
1	A	143	SER	3.1
1	A	6	GLN	3.1
2	B	81	LEU	3.1
1	A	214	PHE	3.0
2	B	262	ASP	3.0
2	B	68	LYS	3.0
1	A	1	GLY	3.0
1	A	177	ASP	2.9
1	A	186	ASN	2.9
1	A	193	ASN	2.8
2	B	223	ARG	2.8
2	B	261	ALA	2.8
1	A	215	PRO	2.8
1	A	158	THR	2.7
1	A	194	LYS	2.7
1	A	208	ILE	2.7
1	A	138	LEU	2.7
1	A	144	SER	2.7
1	A	147	SER	2.7
2	B	128	ARG	2.7
1	A	198	ALA	2.7
2	B	122	PHE	2.6
2	B	28	ASN	2.6
1	A	216	SER	2.6
2	B	26	ASP	2.6
2	B	57	VAL	2.6
2	B	201	LEU	2.5
1	A	164	LYS	2.5
1	A	166	SER	2.5
1	A	127	ASN	2.5
2	B	200	ALA	2.5
1	A	128	ILE	2.4
1	A	160	VAL	2.4
2	B	83	LYS	2.4
2	B	31	TYR	2.4
1	A	129	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	141	SER	2.4
1	A	197	PHE	2.4
1	A	178	MET	2.4
1	A	196	ASP	2.3
2	B	27	MET	2.3
1	A	132	ASP	2.3
2	B	193	GLN	2.3
1	A	185	SER	2.3
1	A	157	GLN	2.3
1	A	205	ASN	2.2
1	A	204	ASN	2.2
2	B	22	LEU	2.2
1	A	172	ASP	2.2
2	B	244	ASP	2.2
1	A	202	ALA	2.2
1	A	179	ARG	2.1
2	B	121	LEU	2.1
1	A	47	GLY	2.1
1	A	140	ASP	2.1
2	B	260	ARG	2.1
1	A	187	SER	2.1
2	B	15	THR	2.1
2	B	23	CYS	2.1
2	B	47	GLY	2.0
1	A	200	ALA	2.0
1	A	146	LYS	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.