



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:15 AM UTC

PDB ID : 8TDX / pdb_00008tdx
Title : TRNM-b.01 in complex with HIV Env fusion peptide
Authors : Olia, A.S.; Kwong, P.D.
Deposited on : 2023-07-05
Resolution : 2.09 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

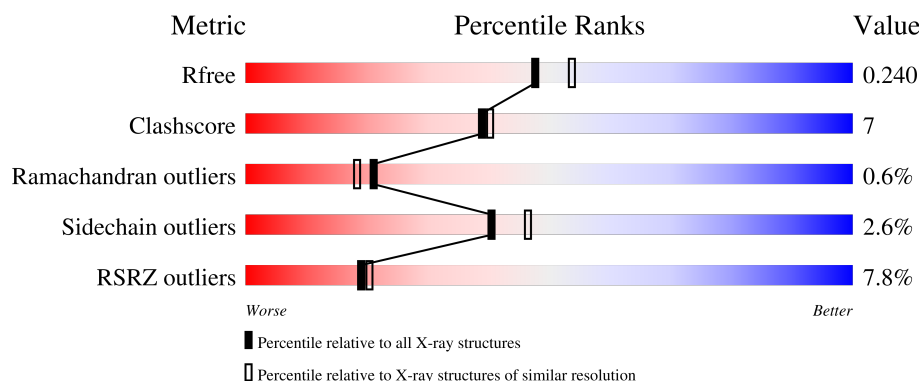
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	227	8% 85% 12% ..
1	B	227	13% 82% 15% ..
2	C	213	3% 82% 18%
2	D	213	5% 80% 18% ..
3	E	8	62% 88% 12%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	F	8	12% 88% 12%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7500 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 TRNM-b.01 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	0	0
			1705	1072	291	335	7			
1	B	223	Total	C	N	O	S	0	0	0
			1705	1072	291	335	7			

- Molecule 2 is a protein called TRNM-b.01 Fab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	213	Total	C	N	O	S	0	0	0
			1620	1015	269	331	5			
2	D	211	Total	C	N	O	S	0	0	0
			1610	1009	267	329	5			

- Molecule 3 is a protein called Env Fusion Peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	8	Total	C	N	O	0	0	0
			51	35	8	8			
3	E	8	Total	C	N	O	0	0	0
			51	35	8	8			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	189	Total	O	0	0
			189	189		
4	C	224	Total	O	0	0
			224	224		
4	D	156	Total	O	0	0
			156	156		
4	B	183	Total	O	0	0
			183	183		

Continued on next page...

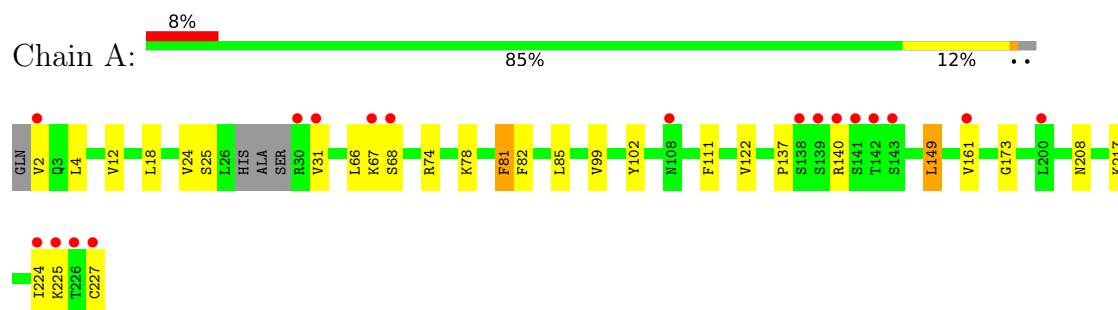
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	3	Total	O	0	0
			3	3		
4	E	3	Total	O	0	0
			3	3		

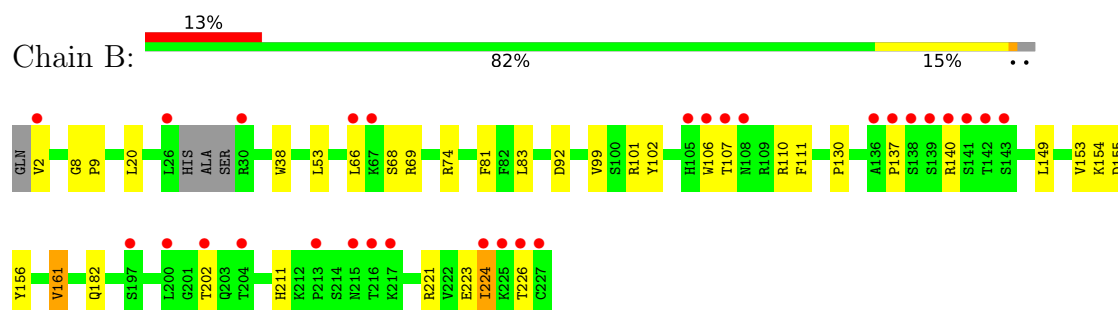
3 Residue-property plots [i](#)

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.

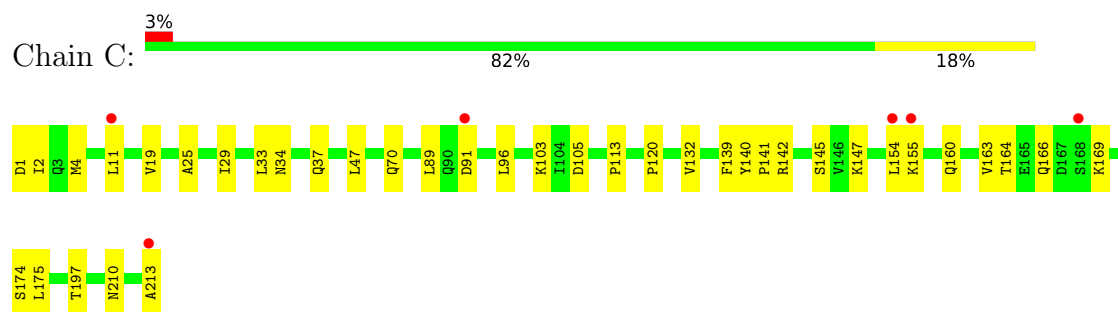
• Molecule 1: TRNM-b.01 Fab Heavy Chain



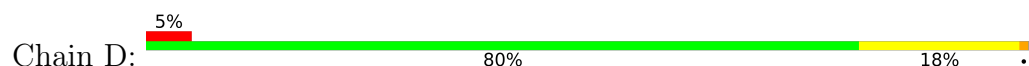
• Molecule 1: TRNM-b.01 Fab Heavy Chain

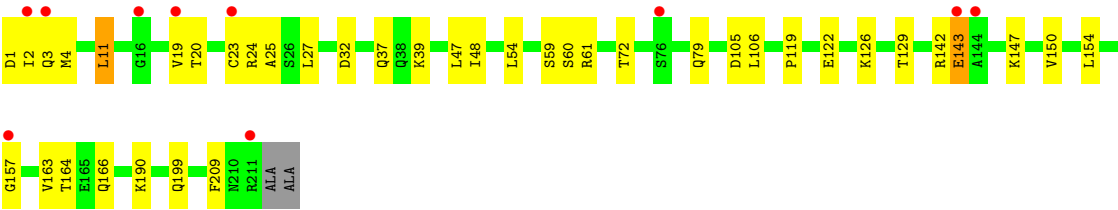


• Molecule 2: TRNM-b.01 Fab Light Chain

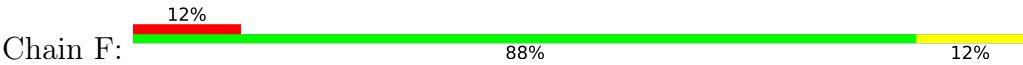


• Molecule 2: TRNM-b.01 Fab Light Chain

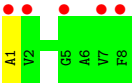
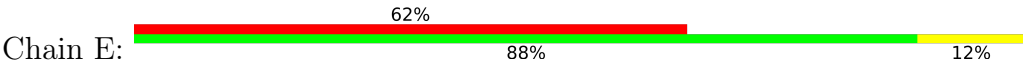




● Molecule 3: Env Fusion Peptide



● Molecule 3: Env Fusion Peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.46Å 88.93Å 129.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.46 – 2.09 44.46 – 2.09	Depositor EDS
% Data completeness (in resolution range)	98.3 (44.46-2.09) 94.5 (44.46-2.09)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.82 (at 2.08Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.204 , 0.240 0.205 , 0.240	Depositor DCC
R_{free} test set	1997 reflections (3.49%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.282	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7500	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 17.76% 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	A	0.17	0/1747	0.41	0/2383
1	B	0.17	0/1747	0.39	0/2383
2	C	0.16	0/1653	0.40	0/2248
2	D	0.19	0/1643	0.44	0/2234
3	E	0.19	0/51	0.44	0/68
3	F	0.28	0/51	0.40	0/68
All	All	0.18	0/6892	0.41	0/9384

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	66	LEU	Peptide

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	1705	0	1653	17	0
1	B	1705	0	1653	25	0
2	C	1620	0	1583	27	0
2	D	1610	0	1573	26	0
3	E	51	0	56	1	0
3	F	51	0	56	0	0
4	A	189	0	0	5	0
4	B	183	0	0	6	1
4	C	224	0	0	7	1
4	D	156	0	0	7	0
4	E	3	0	0	0	0
4	F	3	0	0	0	0
All	All	7500	0	6574	91	1

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 7.

All (91) 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:B:140:ARG:NH2	4:B:301:HOH:O	1.93	1.00
2:C:155:LYS:NZ	4:C:302:HOH:O	2.04	0.84
1:B:155:ASP:OD1	4:B:302:HOH:O	1.99	0.81
1:A:227:CYS:O	4:A:301:HOH:O	2.01	0.77
2:D:157:GLY:O	4:D:301:HOH:O	2.01	0.77
1:B:221:ARG:HH11	1:B:223:GLU:HB3	1.52	0.75
2:C:160:GLN:HE22	1:B:182:GLN:HE22	1.36	0.73
1:A:78:LYS:NZ	4:A:302:HOH:O	2.24	0.71
1:B:223:GLU:O	4:B:303:HOH:O	2.14	0.65
2:D:39:LYS:NZ	4:D:304:HOH:O	2.30	0.64
2:D:37:GLN:HB2	2:D:47:LEU:HD11	1.79	0.64
2:C:11:LEU:HD11	2:C:19:VAL:HG13	1.80	0.63
2:C:147:LYS:NZ	4:C:307:HOH:O	2.29	0.60
1:B:137:PRO:HG3	1:B:149:LEU:HB3	1.82	0.60
2:D:4:MET:HE1	2:D:25:ALA:HB2	1.86	0.58
2:C:1:ASP:N	4:C:301:HOH:O	1.96	0.57
2:C:160:GLN:HE22	1:B:182:GLN:NE2	2.00	0.57
2:D:2:ILE:HG12	2:D:27:LEU:HD22	1.89	0.55
2:D:142:ARG:NH1	2:D:163:VAL:HG21	2.21	0.55
2:D:20:THR:HG23	2:D:72:THR:HG23	1.89	0.55
2:C:145:SER:OG	2:C:197:THR:HB	2.08	0.53
1:B:224:ILE:HD13	1:B:226:THR:HG23	1.90	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:4:MET:HB3	2:D:23:CYS:SG	2.48	0.53
1:B:221:ARG:NH1	1:B:223:GLU:HB3	2.20	0.53
2:D:61:ARG:HH12	2:D:79:GLN:HG2	1.74	0.53
2:C:70:GLN:NE2	4:C:311:HOH:O	2.41	0.53
2:D:39:LYS:HE2	4:D:431:HOH:O	2.08	0.52
1:A:18:LEU:HB3	1:A:85:LEU:HB3	1.91	0.51
2:C:164:THR:HG22	2:C:174:SER:H	1.75	0.51
1:A:224:ILE:H	1:A:224:ILE:HD12	1.75	0.51
2:C:4:MET:HE1	2:C:25:ALA:HB2	1.93	0.51
2:C:163:VAL:HG22	2:C:175:LEU:HD12	1.93	0.51
2:C:210:ASN:HB3	2:C:213:ALA:HB3	1.92	0.51
1:A:208:ASN:HB3	1:A:217:LYS:HE3	1.93	0.51
2:C:103:LYS:NZ	4:C:308:HOH:O	2.34	0.50
2:C:2:ILE:HD13	2:C:29:ILE:HG22	1.93	0.49
1:A:137:PRO:HG3	1:A:149:LEU:HB3	1.94	0.49
2:D:190:LYS:NZ	4:D:302:HOH:O	2.14	0.49
1:A:12:VAL:HG23	1:A:122:VAL:HG22	1.94	0.48
1:A:225:LYS:NZ	2:D:122:GLU:HG3	2.29	0.47
4:C:309:HOH:O	1:B:154:LYS:NZ	2.47	0.47
2:D:24:ARG:NE	4:D:307:HOH:O	2.36	0.47
2:C:142:ARG:NH2	2:C:163:VAL:HG21	2.30	0.47
1:B:106:TRP:HD1	1:B:107:THR:HG23	1.80	0.47
1:A:173:GLY:HA2	4:D:400:HOH:O	2.15	0.46
2:C:89:LEU:HD11	2:C:96:LEU:HB3	1.98	0.46
2:C:140:TYR:CG	2:C:141:PRO:HA	2.50	0.46
1:A:2:VAL:HG11	4:A:446:HOH:O	2.16	0.46
2:D:48:ILE:HD13	2:D:54:LEU:HA	1.97	0.46
2:D:143:GLU:OE1	2:D:199:GLN:HB2	2.15	0.46
1:A:4:LEU:HD22	1:A:24:VAL:HG22	1.98	0.45
2:C:91:ASP:OD2	1:B:101:ARG:NH2	2.48	0.45
1:B:2:VAL:HG21	1:B:102:TYR:OH	2.16	0.45
1:A:2:VAL:HG13	1:A:25:SER:O	2.16	0.45
1:A:102:TYR:HE1	4:A:446:HOH:O	1.98	0.45
1:B:9:PRO:HG2	4:B:391:HOH:O	2.16	0.45
2:D:11:LEU:HD21	2:D:19:VAL:HG13	1.99	0.45
1:B:38:TRP:CD1	1:B:83:LEU:HB2	2.52	0.45
1:B:140:ARG:NH2	4:B:311:HOH:O	2.50	0.45
2:C:37:GLN:HB2	2:C:47:LEU:HD11	1.97	0.44
1:B:110:ARG:O	1:B:110:ARG:HG3	2.14	0.44
1:B:130:PRO:HB3	1:B:156:TYR:HB3	2.00	0.44
1:B:8:GLY:HA3	1:B:20:LEU:HD23	1.99	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:105:ASP:HB2	2:C:166:GLN:OE1	2.18	0.44
2:D:142:ARG:HH11	2:D:163:VAL:HG21	1.81	0.44
2:D:147:LYS:HE3	2:D:154:LEU:HD13	2.00	0.44
2:C:145:SER:HG	2:C:197:THR:HB	1.83	0.44
2:D:61:ARG:HH12	2:D:79:GLN:CG	2.31	0.43
1:B:74:ARG:HA	1:B:81:PHE:HA	1.99	0.43
1:B:153:VAL:HG11	1:B:161:VAL:HG11	1.99	0.43
1:B:161:VAL:HG22	1:B:211:HIS:HB2	1.99	0.43
2:C:169:LYS:HG3	4:C:424:HOH:O	2.19	0.43
2:D:163:VAL:HG22	2:D:164:THR:O	2.17	0.43
1:A:74:ARG:HA	1:A:81:PHE:HA	1.99	0.42
2:D:32:ASP:OD2	3:E:1:ALA:HB2	2.19	0.42
2:C:120:PRO:HD3	2:C:132:VAL:HG22	2.02	0.42
2:D:61:ARG:H	2:D:61:ARG:HG3	1.64	0.42
2:C:33:LEU:HG	2:C:34:ASN:N	2.34	0.42
2:D:59:SER:OG	2:D:61:ARG:HG3	2.19	0.42
2:D:105:ASP:HB2	2:D:166:GLN:OE1	2.19	0.42
2:C:164:THR:CG2	2:C:174:SER:H	2.32	0.42
2:C:147:LYS:HD2	2:C:154:LEU:HD13	2.02	0.42
1:A:137:PRO:O	1:A:224:ILE:HG23	2.20	0.41
1:B:154:LYS:HE3	4:B:386:HOH:O	2.20	0.41
1:A:82:PHE:HB3	4:A:425:HOH:O	2.20	0.41
2:D:119:PRO:HB3	2:D:209:PHE:CE2	2.56	0.41
1:B:69:ARG:NH2	1:B:92:ASP:OD2	2.49	0.41
2:D:142:ARG:NH2	4:D:314:HOH:O	2.54	0.41
2:C:113:PRO:HB3	2:C:139:PHE:HB3	2.03	0.40
1:A:66:LEU:O	1:A:68:SER:N	2.55	0.40
1:B:102:TYR:O	1:B:110:ARG:O	2.39	0.40

All (1) 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)
4:C:489:HOH:O	4:B:462:HOH:O[3_554]	2.00	0.20

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	219/227 (96%)	213 (97%)	3 (1%)	3 (1%)	9	5
1	B	219/227 (96%)	209 (95%)	9 (4%)	1 (0%)	24	22
2	C	211/213 (99%)	206 (98%)	5 (2%)	0	100	100
2	D	209/213 (98%)	200 (96%)	8 (4%)	1 (0%)	24	22
3	E	6/8 (75%)	6 (100%)	0	0	100	100
3	F	6/8 (75%)	6 (100%)	0	0	100	100
All	All	870/896 (97%)	840 (97%)	25 (3%)	5 (1%)	21	18

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	111	PHE
1	A	67	LYS
1	A	111	PHE
2	D	143	GLU
1	A	31	VAL

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	197/203 (97%)	192 (98%)	5 (2%)	42	48
1	B	197/203 (97%)	191 (97%)	6 (3%)	36	41
2	C	186/186 (100%)	186 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	186/186 (100%)	178 (96%)	8 (4%)	26	27
3	E	4/4 (100%)	4 (100%)	0	100	100
3	F	4/4 (100%)	3 (75%)	1 (25%)	0	0
All	All	774/786 (98%)	754 (97%)	20 (3%)	40	46

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	PHE
1	A	99	VAL
1	A	140	ARG
1	A	149	LEU
1	A	161	VAL
2	D	1	ASP
2	D	3	GLN
2	D	11	LEU
2	D	60	SER
2	D	106	LEU
2	D	126	LYS
2	D	129	THR
2	D	150	VAL
1	B	53	LEU
1	B	68	SER
1	B	99	VAL
1	B	161	VAL
1	B	202	THR
1	B	224	ILE
3	F	2	VAL

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

Mol	Chain	Res	Type
2	C	70	GLN
2	C	124	GLN
2	C	137	ASN
2	C	138	ASN
2	C	160	GLN
2	D	30	ASN
1	B	3	GLN
1	B	5	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	119	GLN
1	B	211	HIS

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	223/227 (98%)	0.29	18 (8%) 18 19	17, 28, 57, 95	0
1	B	223/227 (98%)	0.59	29 (13%) 7 7	17, 30, 66, 94	0
2	C	213/213 (100%)	0.09	6 (2%) 55 57	18, 26, 39, 63	0
2	D	211/213 (99%)	0.50	10 (4%) 36 38	19, 32, 46, 58	0
3	E	8/8 (100%)	3.06	5 (62%) 0 0	31, 42, 66, 69	0
3	F	8/8 (100%)	0.36	1 (12%) 8 8	21, 25, 30, 31	0
All	All	886/896 (98%)	0.39	69 (7%) 19 20	17, 29, 53, 95	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	1	ALA	8.5
1	B	106	TRP	6.9
1	B	227	CYS	6.1
1	A	141	SER	5.7
1	B	226	THR	5.2
2	D	143	GLU	5.0
1	B	141	SER	4.8
1	B	67	LYS	4.8
1	A	2	VAL	4.8
3	E	8	PHE	4.7
1	A	140	ARG	4.2
1	B	138	SER	4.2
1	A	226	THR	3.9
1	B	2	VAL	3.9
1	B	108	ASN	3.9
1	B	142	THR	3.9
2	C	155	LYS	3.8
1	A	142	THR	3.8
1	B	107	THR	3.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	227	CYS	3.7
1	A	67	LYS	3.6
1	B	66	LEU	3.5
2	D	2	ILE	3.4
1	B	26	LEU	3.4
1	B	213	PRO	3.4
1	B	139	SER	3.3
2	D	157	GLY	3.3
1	A	225	LYS	3.2
1	B	204	THR	3.1
1	A	143	SER	3.1
3	E	7	VAL	3.1
3	E	2	VAL	3.0
1	B	224	ILE	2.8
2	D	144	ALA	2.8
1	B	137	PRO	2.8
1	B	216	THR	2.8
1	B	143	SER	2.7
3	E	5	GLY	2.7
1	A	68	SER	2.7
1	A	139	SER	2.7
2	D	23	CYS	2.7
2	C	154	LEU	2.6
1	A	224	ILE	2.5
1	B	215	ASN	2.5
1	B	105	HIS	2.5
2	C	213	ALA	2.5
1	B	30	ARG	2.4
1	A	31	VAL	2.4
2	D	76	SER	2.4
1	B	200	LEU	2.4
1	B	140	ARG	2.4
1	B	136	ALA	2.3
2	D	3	GLN	2.3
1	B	197	SER	2.3
1	B	202	THR	2.3
1	A	108	ASN	2.2
2	D	211	ARG	2.2
1	B	217	LYS	2.2
3	F	1	ALA	2.2
2	C	91	ASP	2.2
2	D	19	VAL	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	C	11	LEU	2.1
1	A	138	SER	2.1
2	C	168	SER	2.1
1	A	200	LEU	2.1
1	A	30	ARG	2.0
2	D	16	GLY	2.0
1	A	161	VAL	2.0
1	B	225	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.