



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:07 AM UTC

PDB ID : 6WIR / pdb_00006wir
Title : Fab antigen complex
Authors : Antonysamy, S.
Deposited on : 2020-04-10
Resolution : 2.96 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

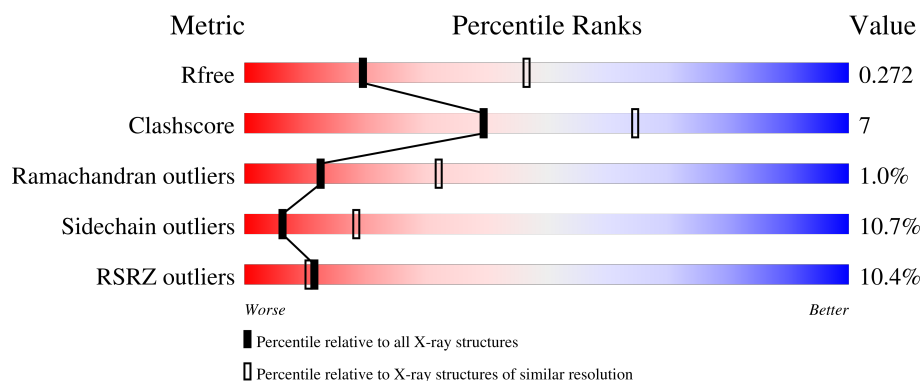
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.96 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	236	6% 69% 24% . .
2	B	213	6% 70% 28% .
3	C	161	18% 43% 11% . 45%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4125 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 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	227	Total	C	N	O	S	0	0	0
			1721	1095	288	331	7			

- Molecule 2 is a protein called Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	0	0
			1610	1002	275	328	5			

- Molecule 3 is a protein called Interleukin-17A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	89	Total	C	N	O	S	0	0	0
			722	452	133	131	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	133	HIS	-	expression tag	UNP Q16552
C	134	HIS	-	expression tag	UNP Q16552
C	135	HIS	-	expression tag	UNP Q16552
C	136	HIS	-	expression tag	UNP Q16552
C	137	HIS	-	expression tag	UNP Q16552
C	138	HIS	-	expression tag	UNP Q16552

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	34	Total	O	0	0
			34	34		
4	B	36	Total	O	0	0
			36	36		

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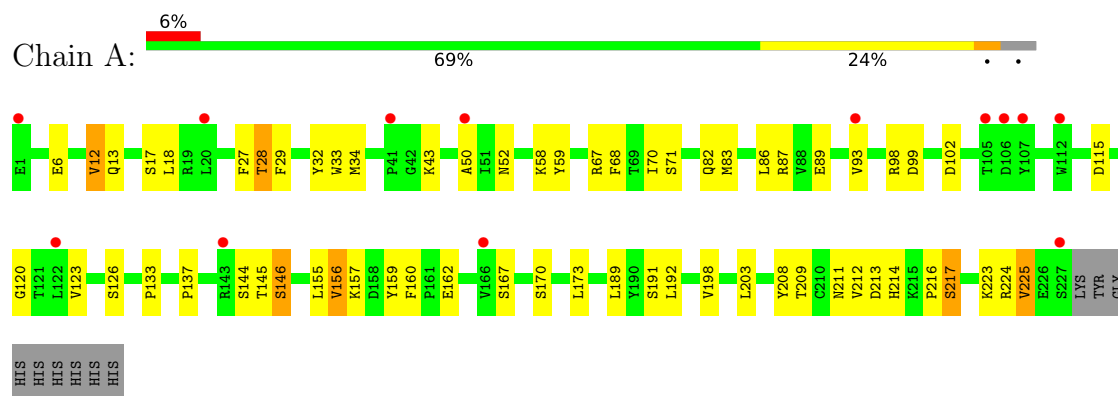
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	2	Total	O	0	0
			2	2		

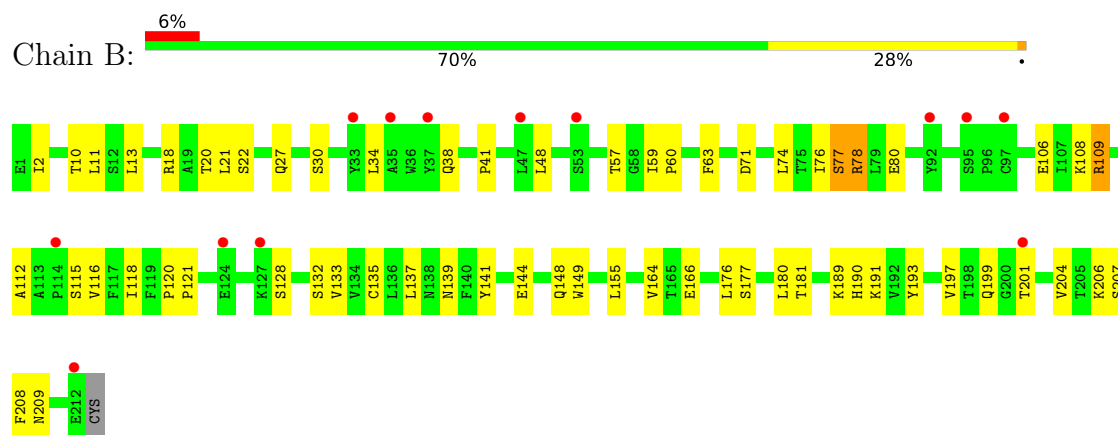
3 Residue-property plots

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.

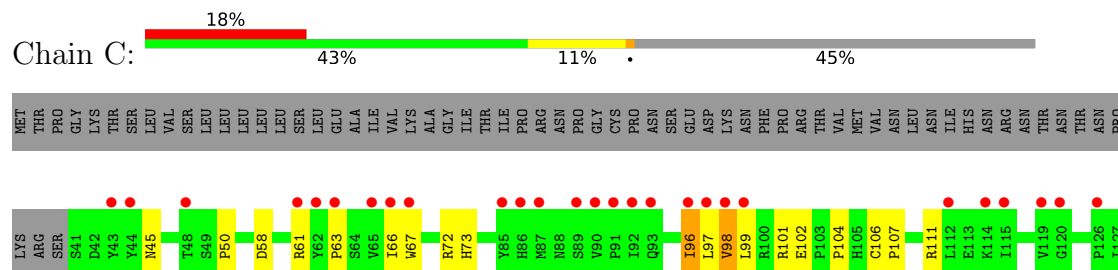
• Molecule 1: Fab heavy chain



• Molecule 2: Fab light chain



• Molecule 3: Interleukin-17A



Y128
H129
HIS
VAL
ALA
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	73.40Å 73.40Å 326.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.61 – 2.96 71.61 – 2.96	Depositor EDS
% Data completeness (in resolution range)	100.0 (71.61-2.96) 100.0 (71.61-2.96)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.96Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.217 , 0.260 0.227 , 0.272	Depositor DCC
R_{free} test set	977 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	94.3	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 135.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4125	wwPDB-VP
Average B, all atoms (Å ²)	111.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% 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	0.92	2/1766 (0.1%)	1.32	10/2411 (0.4%)
2	B	0.89	0/1643	1.30	9/2230 (0.4%)
3	C	0.91	1/743 (0.1%)	1.31	0/1015
All	All	0.91	3/4152 (0.1%)	1.31	19/5656 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	155	LEU	CA-C	6.46	1.61	1.52
1	A	156	VAL	CA-C	5.82	1.60	1.52
3	C	96	ILE	CA-C	5.44	1.57	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	28	THR	CA-C-N	6.67	129.52	120.38
1	A	28	THR	C-N-CA	6.67	129.52	120.38
1	A	102	ASP	CA-CB-CG	6.34	118.94	112.60
2	B	201	THR	CA-C-N	5.76	129.44	120.75
2	B	201	THR	C-N-CA	5.76	129.44	120.75
2	B	77	SER	N-CA-C	5.75	117.22	111.07
1	A	157	LYS	CA-C-N	5.48	129.96	122.34
1	A	157	LYS	C-N-CA	5.48	129.96	122.34
1	A	115	ASP	CA-CB-CG	5.37	117.97	112.60
2	B	76	ILE	CA-C-N	5.24	127.25	120.44
2	B	76	ILE	C-N-CA	5.24	127.25	120.44
2	B	78	ARG	CA-C-N	5.16	128.42	121.05
2	B	78	ARG	C-N-CA	5.16	128.42	121.05
1	A	213	ASP	CA-CB-CG	5.13	117.73	112.60
2	B	180	LEU	N-CA-C	-5.13	102.16	110.32
1	A	29	PHE	CA-CB-CG	5.12	118.92	113.80
1	A	146	SER	CA-C-N	5.09	131.25	121.54
1	A	146	SER	C-N-CA	5.09	131.25	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	144	GLU	CB-CG-CD	5.01	121.12	112.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1721	0	1644	23	0
2	B	1610	0	1550	23	0
3	C	722	0	688	7	0
4	A	34	0	0	1	0
4	B	36	0	0	0	0
4	C	2	0	0	0	0
All	All	4125	0	3882	53	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 7.

All (53) 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:38:GLN:HB2	2:B:48:LEU:HD11	1.60	0.83
2:B:18:ARG:HG3	2:B:77:SER:HA	1.69	0.73
1:A:67:ARG:HH12	1:A:87:ARG:HG3	1.55	0.71
2:B:164:VAL:HG22	2:B:176:LEU:HD12	1.74	0.69
3:C:101:ARG:HD3	3:C:106:CYS:H	1.57	0.68
1:A:12:VAL:HG21	1:A:18:LEU:HG	1.74	0.68
2:B:109:ARG:NH2	2:B:112:ALA:HB2	2.12	0.64
1:A:160:PHE:HB2	1:A:189:LEU:HD23	1.80	0.63
2:B:108:LYS:HA	2:B:141:TYR:OH	2.02	0.60
3:C:63:PRO:HD2	3:C:98:VAL:HG22	1.85	0.58
1:A:50:ALA:HB3	1:A:59:TYR:HB2	1.86	0.57
1:A:214:HIS:CE1	1:A:217:SER:HB3	2.41	0.55
2:B:41:PRO:HG3	2:B:166:GLU:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LYS:HE3	1:A:70:ILE:O	2.07	0.54
1:A:214:HIS:CD2	1:A:216:PRO:HD2	2.42	0.54
2:B:191:LYS:O	2:B:209:ASN:HA	2.07	0.54
2:B:164:VAL:CG2	2:B:176:LEU:HD12	2.36	0.54
2:B:148:GLN:HB3	2:B:155:LEU:HD11	1.90	0.52
1:A:83:MET:HB3	1:A:86:LEU:HD21	1.93	0.51
2:B:21:LEU:HD22	2:B:74:LEU:HD23	1.93	0.51
1:A:156:VAL:HG13	1:A:212:VAL:HG11	1.94	0.49
2:B:109:ARG:HH22	2:B:112:ALA:HB2	1.77	0.49
2:B:2:ILE:HG12	2:B:27:GLN:HB2	1.95	0.48
1:A:32:TYR:HB2	1:A:34:MET:HE2	1.95	0.48
3:C:50:PRO:HB2	3:C:72:ARG:HH11	1.80	0.47
1:A:133:PRO:HB3	1:A:159:TYR:HB3	1.97	0.47
1:A:17:SER:HA	1:A:83:MET:O	2.15	0.46
2:B:116:VAL:HG12	2:B:206:LYS:HG3	1.97	0.46
1:A:83:MET:HE1	1:A:123:VAL:HG21	1.97	0.45
1:A:214:HIS:ND1	1:A:217:SER:HB3	2.30	0.45
1:A:137:PRO:HB2	1:A:225:VAL:HG13	1.98	0.45
3:C:96:ILE:HG13	3:C:97:LEU:H	1.81	0.45
1:A:34:MET:HE1	1:A:98:ARG:HG3	1.99	0.44
2:B:120:PRO:HB3	2:B:208:PHE:CZ	2.52	0.44
2:B:190:HIS:HB2	2:B:193:TYR:OH	2.18	0.44
2:B:48:LEU:HD23	2:B:59:ILE:HD12	2.00	0.43
3:C:50:PRO:HB2	3:C:72:ARG:HD2	2.01	0.43
3:C:58:ASP:HB3	3:C:61:ARG:HB2	1.99	0.43
1:A:198:VAL:HG11	1:A:208:TYR:CE1	2.55	0.42
1:A:6:GLU:CD	1:A:120:GLY:H	2.27	0.42
1:A:67:ARG:NH1	1:A:87:ARG:HG3	2.29	0.42
2:B:60:PRO:HG2	2:B:63:PHE:HD1	1.83	0.42
3:C:101:ARG:HG3	3:C:104:PRO:HA	2.02	0.42
2:B:11:LEU:HG	2:B:13:LEU:HD13	2.02	0.42
2:B:137:LEU:HD11	2:B:197:VAL:CG2	2.50	0.42
1:A:68:PHE:HA	1:A:82:GLN:O	2.21	0.41
2:B:121:PRO:HB3	2:B:132:SER:H	1.85	0.41
2:B:137:LEU:HD22	2:B:176:LEU:HD22	2.03	0.41
2:B:118:ILE:HG22	2:B:206:LYS:HB3	2.03	0.41
2:B:135:CYS:HB2	2:B:149:TRP:CH2	2.56	0.41
1:A:27:PHE:HB2	4:A:317:HOH:O	2.21	0.41
1:A:209:THR:HA	1:A:223:LYS:O	2.21	0.41
1:A:33:TRP:HB2	1:A:99:ASP:HB2	2.03	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	225/236 (95%)	201 (89%)	21 (9%)	3 (1%)	9	26
2	B	210/213 (99%)	196 (93%)	13 (6%)	1 (0%)	24	49
3	C	87/161 (54%)	76 (87%)	10 (12%)	1 (1%)	11	29
All	All	522/610 (86%)	473 (91%)	44 (8%)	5 (1%)	12	32

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	146	SER
2	B	139	ASN
1	A	145	THR
1	A	144	SER
3	C	107	PRO

5.3.2 Protein sidechains [i](#)

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	186/203 (92%)	166 (89%)	20 (11%)	6	18
2	B	179/184 (97%)	159 (89%)	20 (11%)	6	17
3	C	83/150 (55%)	75 (90%)	8 (10%)	8	22
All	All	448/537 (83%)	400 (89%)	48 (11%)	6	18

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	13	GLN
1	A	28	THR
1	A	43	LYS
1	A	52	ASN
1	A	71	SER
1	A	89	GLU
1	A	93	VAL
1	A	126	SER
1	A	162	GLU
1	A	167	SER
1	A	170	SER
1	A	173	LEU
1	A	191	SER
1	A	192	LEU
1	A	203	LEU
1	A	211	ASN
1	A	217	SER
1	A	224	ARG
1	A	225	VAL
2	B	10	THR
2	B	20	THR
2	B	22	SER
2	B	30	SER
2	B	34	LEU
2	B	57	THR
2	B	71	ASP
2	B	78	ARG
2	B	80	GLU
2	B	106	GLU
2	B	109	ARG
2	B	115	SER
2	B	128	SER
2	B	133	VAL
2	B	177	SER
2	B	181	THR
2	B	189	LYS
2	B	199	GLN
2	B	204	VAL
2	B	207	SER
3	C	45	ASN
3	C	66	ILE
3	C	67	TRP

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Mol	Chain	Res	Type
3	C	73	HIS
3	C	98	VAL
3	C	99	LEU
3	C	102	GLU
3	C	111	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	ASN
2	B	27	GLN
2	B	39	GLN
2	B	138	ASN
2	B	161	GLN
2	B	199	GLN
2	B	209	ASN
3	C	45	ASN
3	C	86	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 ⓘ

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	227/236 (96%)	0.43	13 (5%)	29 24	68, 99, 130, 145	0
2	B	212/213 (99%)	0.37	13 (6%)	27 22	71, 100, 125, 145	0
3	C	89/161 (55%)	1.61	29 (32%)	1 1	97, 132, 244, 259	0
All	All	528/610 (86%)	0.61	55 (10%)	11 10	68, 103, 190, 259	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	227	SER	6.7
3	C	93	GLN	5.9
2	B	201	THR	5.2
3	C	97	LEU	5.2
3	C	99	LEU	5.0
3	C	90	VAL	4.7
3	C	63	PRO	4.4
3	C	112	LEU	4.0
1	A	112	TRP	3.8
3	C	128	VAL	3.8
1	A	41	PRO	3.8
1	A	1	GLU	3.6
2	B	97	CYS	3.4
3	C	96	ILE	3.4
2	B	35	ALA	3.3
3	C	65	VAL	3.2
2	B	124	GLU	3.2
3	C	66	ILE	3.2
3	C	92	ILE	3.2
3	C	43	TYR	3.1
1	A	50	ALA	3.1
2	B	212	GLU	2.9
3	C	119	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	114	PRO	2.7
3	C	67	TRP	2.7
3	C	44	TYR	2.7
1	A	122	LEU	2.7
3	C	61	ARG	2.7
3	C	91	PRO	2.6
3	C	120	GLY	2.6
3	C	126	PRO	2.6
3	C	89	SER	2.5
3	C	48	THR	2.5
2	B	53	SER	2.4
3	C	114	LYS	2.4
1	A	107	TYR	2.4
3	C	98	VAL	2.4
3	C	129	HIS	2.4
3	C	62	TYR	2.3
1	A	105	THR	2.3
1	A	166	VAL	2.3
2	B	37	TYR	2.3
1	A	143	ARG	2.2
2	B	92	TYR	2.2
2	B	95	SER	2.2
3	C	86	HIS	2.2
2	B	127	LYS	2.1
1	A	106	ASP	2.1
3	C	85	TYR	2.1
1	A	93	VAL	2.1
2	B	33	TYR	2.1
3	C	115	ILE	2.0
1	A	20	LEU	2.0
2	B	47	LEU	2.0
3	C	87	MET	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.