



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2022 – 10:02 AM EDT

PDB ID : 7T13  
Title : Hexameric HIV-1 (M-group) CA Q50Y mutant  
Authors : Jacques, D.A.; James, L.C.  
Deposited on : 2021-12-01  
Resolution : 3.15 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

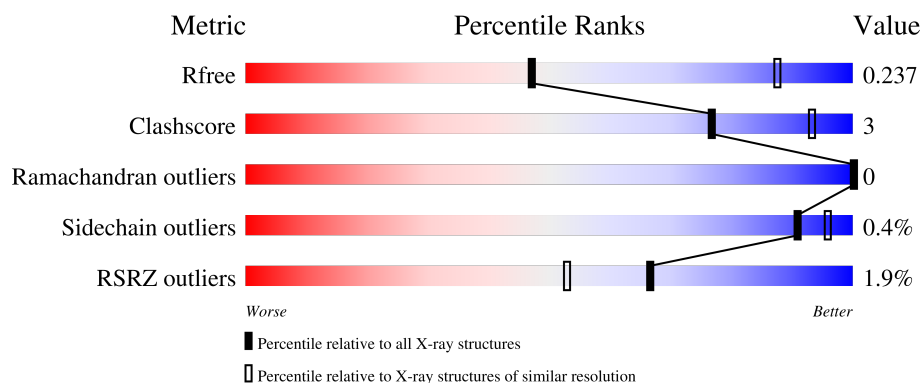
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.15 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.



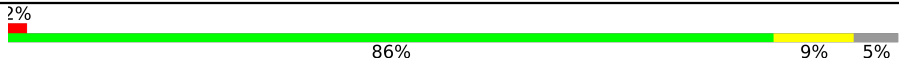
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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  $\geq 3$ , 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	231	
1	B	231	
1	C	231	
1	D	231	
1	E	231	

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Mol	Chain	Length	Quality of chain
1	F	231	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '2%', a large green segment labeled '86%', a yellow segment labeled '9%', and a small grey segment at the end labeled '5%'.

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 9907 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 Capsid protein p24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	9	0
			1695	1070	294	316	15			
1	B	211	Total	C	N	O	S	0	8	0
			1632	1029	281	307	15			
1	C	210	Total	C	N	O	S	0	8	0
			1624	1025	279	305	15			
1	D	211	Total	C	N	O	S	0	9	0
			1634	1032	281	306	15			
1	E	211	Total	C	N	O	S	0	9	0
			1633	1032	280	306	15			
1	F	219	Total	C	N	O	S	0	9	0
			1689	1068	291	315	15			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14	CYS	ALA	engineered mutation	UNP B6DRA0
A	45	CYS	GLU	engineered mutation	UNP B6DRA0
A	50	TYR	GLN	engineered mutation	UNP B6DRA0
A	184	ALA	TRP	engineered mutation	UNP B6DRA0
A	185	ALA	MET	engineered mutation	UNP B6DRA0
B	14	CYS	ALA	engineered mutation	UNP B6DRA0
B	45	CYS	GLU	engineered mutation	UNP B6DRA0
B	50	TYR	GLN	engineered mutation	UNP B6DRA0
B	184	ALA	TRP	engineered mutation	UNP B6DRA0
B	185	ALA	MET	engineered mutation	UNP B6DRA0
C	14	CYS	ALA	engineered mutation	UNP B6DRA0
C	45	CYS	GLU	engineered mutation	UNP B6DRA0
C	50	TYR	GLN	engineered mutation	UNP B6DRA0
C	184	ALA	TRP	engineered mutation	UNP B6DRA0
C	185	ALA	MET	engineered mutation	UNP B6DRA0
D	14	CYS	ALA	engineered mutation	UNP B6DRA0
D	45	CYS	GLU	engineered mutation	UNP B6DRA0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	50	TYR	GLN	engineered mutation	UNP B6DRA0
D	184	ALA	TRP	engineered mutation	UNP B6DRA0
D	185	ALA	MET	engineered mutation	UNP B6DRA0
E	14	CYS	ALA	engineered mutation	UNP B6DRA0
E	45	CYS	GLU	engineered mutation	UNP B6DRA0
E	50	TYR	GLN	engineered mutation	UNP B6DRA0
E	184	ALA	TRP	engineered mutation	UNP B6DRA0
E	185	ALA	MET	engineered mutation	UNP B6DRA0
F	14	CYS	ALA	engineered mutation	UNP B6DRA0
F	45	CYS	GLU	engineered mutation	UNP B6DRA0
F	50	TYR	GLN	engineered mutation	UNP B6DRA0
F	184	ALA	TRP	engineered mutation	UNP B6DRA0
F	185	ALA	MET	engineered mutation	UNP B6DRA0

### 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.

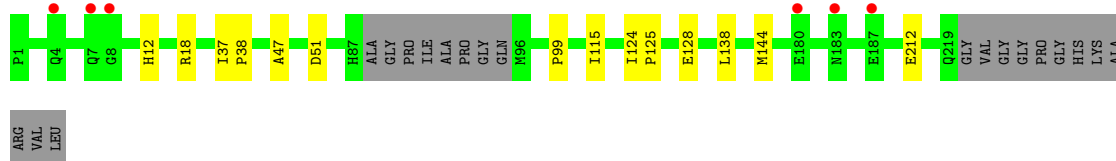
- Molecule 1: Capsid protein p24

Chain A: 




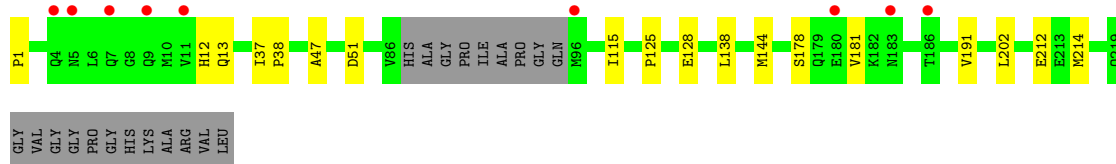
- Molecule 1: Capsid protein p24

Chain B: 




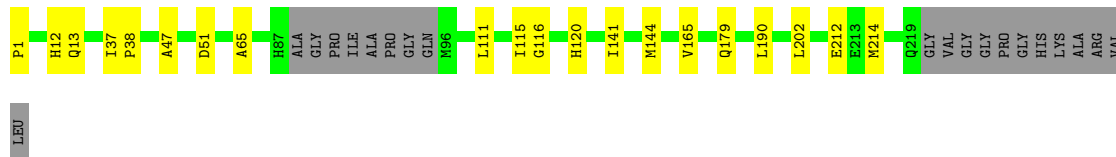
- Molecule 1: Capsid protein p24

Chain C: 

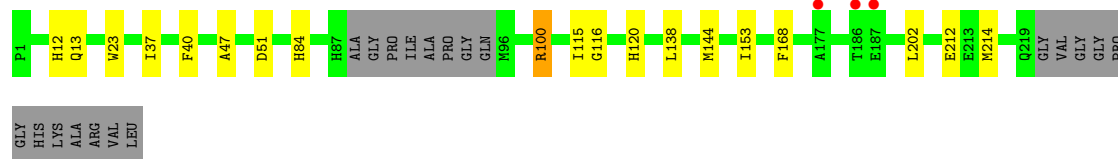
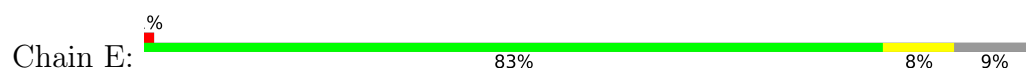


- Molecule 1: Capsid protein p24

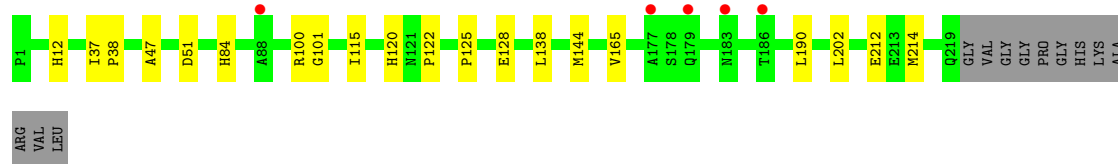
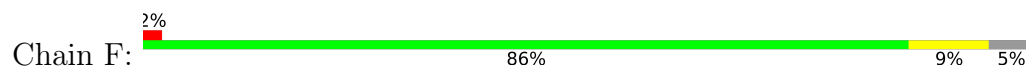
Chain D: 



- Molecule 1: Capsid protein p24



- Molecule 1: Capsid protein p24



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.29Å 156.61Å 118.89Å 90.00° 107.33° 90.00°	Depositor
Resolution (Å)	78.30 – 3.15 78.30 – 3.15	Depositor EDS
% Data completeness (in resolution range)	94.8 (78.30-3.15) 94.9 (78.30-3.15)	Depositor EDS
$R_{merge}$	0.20	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.61 (at 3.13Å)	Xtriage
Refinement program	PHENIX 1.14	Depositor
R, $R_{free}$	0.205 , 0.239 0.204 , 0.237	Depositor DCC
$R_{free}$ test set	1231 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	42.7	Xtriage
Anisotropy	0.447	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 33.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	9907	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.25	0/1739	0.40	0/2368
1	B	0.25	0/1669	0.39	0/2269
1	C	0.25	0/1661	0.39	0/2258
1	D	0.24	0/1674	0.39	0/2277
1	E	0.25	0/1673	0.39	0/2275
1	F	0.25	0/1733	0.40	0/2360
All	All	0.25	0/10149	0.39	0/13807

There are no bond length outliers.

There are no bond angle outliers.

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	1695	0	1681	11	0
1	B	1632	0	1602	9	0
1	C	1624	0	1597	11	0
1	D	1634	0	1612	11	0
1	E	1633	0	1606	11	0
1	F	1689	0	1669	14	0
All	All	9907	0	9767	60	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 (60) 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:E:84:HIS:O	1:E:100:ARG:NH2	2.23	0.70
1:F:84:HIS:O	1:F:100:ARG:NH2	2.30	0.64
1:A:99:PRO:HG3	1:A:124:ILE:HG21	1.80	0.64
1:C:144:MET:HE1	1:D:212:GLU:HG3	1.82	0.62
1:E:144:MET:HE1	1:F:212:GLU:HG3	1.82	0.62
1:F:12:HIS:HB2	1:F:115:ILE:HD11	1.83	0.59
1:B:144:MET:HE1	1:C:212:GLU:HG3	1.85	0.59
1:B:12:HIS:HB2	1:B:115:ILE:HD11	1.83	0.59
1:C:178:SER:HB3	1:C:181:VAL:HG23	1.84	0.59
1:C:12:HIS:HB2	1:C:115:ILE:HD11	1.86	0.57
1:A:144:MET:HE1	1:B:212:GLU:HG3	1.89	0.55
1:F:37:ILE:HD12	1:F:138:LEU:HB3	1.90	0.53
1:F:47:ALA:HB1	1:F:51:ASP:HB2	1.90	0.53
1:D:12:HIS:HB2	1:D:115:ILE:HD11	1.90	0.53
1:E:12:HIS:HB2	1:E:115:ILE:HD11	1.90	0.52
1:A:3:VAL:HG22	1:A:13:GLN:HB2	1.91	0.52
1:E:47:ALA:HB1	1:E:51:ASP:HB2	1.90	0.52
1:A:96:MET:HE1	1:A:116:GLY:HA3	1.92	0.52
1:B:99:PRO:HG3	1:B:124:ILE:HG21	1.92	0.51
1:D:37:ILE:HB	1:D:38[A]:PRO:HD3	1.92	0.51
1:C:125:PRO:HB2	1:C:128:GLU:HB2	1.93	0.50
1:A:12:HIS:HB2	1:A:115:ILE:HD11	1.94	0.50
1:A:212:GLU:HG3	1:F:144:MET:HE1	1.92	0.50
1:A:96:MET:HG3	1:A:97:ARG:H	1.77	0.50
1:D:144:MET:HE1	1:E:212:GLU:HG3	1.94	0.49
1:D:1:PRO:HB2	1:D:13:GLN:HE21	1.76	0.49
1:C:37:ILE:HB	1:C:38[A]:PRO:HD3	1.95	0.49
1:D:202:LEU:HD22	1:D:214[A]:MET:HE3	1.94	0.48
1:F:125:PRO:HB2	1:F:128:GLU:HB2	1.94	0.48
1:B:47:ALA:HB1	1:B:51:ASP:HB2	1.96	0.47
1:A:81:ASP:HA	1:A:100:ARG:HH12	1.80	0.46
1:D:165:VAL:HG22	1:D:190:LEU:HD11	1.96	0.46
1:E:100:ARG:HA	1:E:100:ARG:HE	1.81	0.46
1:A:37:ILE:HD12	1:A:138:LEU:HB3	1.99	0.45
1:F:120:HIS:ND1	1:F:122:PRO:O	2.47	0.45
1:E:37:ILE:HD12	1:E:138:LEU:HB3	1.99	0.45
1:D:65:ALA:HB1	1:D:141:ILE:HD13	1.99	0.44
1:E:116:GLY:O	1:E:120:HIS:HB2	2.17	0.44
1:D:12:HIS:CG	1:D:111:LEU:HD21	2.52	0.44
1:B:37:ILE:HB	1:B:38[A]:PRO:HD3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:191:VAL:HG22	1:C:202:LEU:HD13	2.00	0.43
1:B:37:ILE:HD12	1:B:138:LEU:HB3	2.01	0.43
1:C:202:LEU:HD22	1:C:214[A]:MET:HE3	2.01	0.42
1:A:18:ARG:HG2	1:B:18:ARG:NH2	2.34	0.42
1:B:125:PRO:HB2	1:B:128:GLU:HB2	2.01	0.42
1:D:47:ALA:HB1	1:D:51:ASP:HB2	2.01	0.42
1:D:116:GLY:O	1:D:120:HIS:HB2	2.19	0.42
1:C:1:PRO:HB2	1:C:13:GLN:HE21	1.83	0.42
1:F:37:ILE:HB	1:F:38[A]:PRO:HD3	2.02	0.42
1:E:202:LEU:HD22	1:E:214[A]:MET:HE3	2.02	0.42
1:C:37:ILE:HD12	1:C:138:LEU:HB3	2.01	0.41
1:E:153:ILE:HG21	1:E:168:PHE:HA	2.01	0.41
1:A:47:ALA:HB1	1:A:51:ASP:HB2	2.02	0.41
1:F:100:ARG:HG3	1:F:101:GLY:H	1.85	0.41
1:F:165:VAL:HG22	1:F:190:LEU:HD11	2.01	0.41
1:F:202:LEU:HD22	1:F:214[A]:MET:HE3	2.03	0.41
1:C:47:ALA:HB1	1:C:51:ASP:HB2	2.02	0.41
1:F:100:ARG:HG3	1:F:101:GLY:N	2.36	0.41
1:F:100:ARG:HE	1:F:100:ARG:HA	1.86	0.41
1:E:23:TRP:CZ3	1:E:40:PHE:HB2	2.56	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	219/231 (95%)	213 (97%)	6 (3%)	0	100	100
1	B	208/231 (90%)	200 (96%)	8 (4%)	0	100	100
1	C	207/231 (90%)	199 (96%)	8 (4%)	0	100	100
1	D	209/231 (90%)	202 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	209/231 (90%)	202 (97%)	7 (3%)	0	100	100
1	F	219/231 (95%)	211 (96%)	8 (4%)	0	100	100
All	All	1271/1386 (92%)	1227 (96%)	44 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	184/193 (95%)	183 (100%)	1 (0%)	88	95
1	B	176/193 (91%)	176 (100%)	0	100	100
1	C	175/193 (91%)	175 (100%)	0	100	100
1	D	177/193 (92%)	176 (99%)	1 (1%)	86	94
1	E	176/193 (91%)	174 (99%)	2 (1%)	73	88
1	F	182/193 (94%)	182 (100%)	0	100	100
All	All	1070/1158 (92%)	1066 (100%)	4 (0%)	91	96

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

Mol	Chain	Res	Type
1	A	143	ARG
1	D	179	GLN
1	E	13	GLN
1	E	100	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	13	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 monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	219/231 (94%)	-0.27	1 (0%) 91 86	15, 32, 70, 176	0
1	B	211/231 (91%)	-0.19	6 (2%) 53 36	14, 33, 74, 90	1 (0%)
1	C	210/231 (90%)	-0.15	9 (4%) 35 21	16, 34, 82, 119	0
1	D	211/231 (91%)	-0.30	0 100 100	18, 34, 73, 94	0
1	E	211/231 (91%)	-0.31	3 (1%) 75 63	16, 33, 70, 87	0
1	F	219/231 (94%)	-0.21	5 (2%) 60 46	14, 31, 71, 96	0
All	All	1281/1386 (92%)	-0.24	24 (1%) 66 53	14, 33, 74, 176	1 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	177	ALA	4.6
1	C	4	GLN	4.1
1	C	96	MET	3.3
1	C	7	GLN	3.2
1	F	88	ALA	3.1
1	B	187	GLU	2.9
1	B	180	GLU	2.7
1	F	177	ALA	2.6
1	F	186	THR	2.6
1	C	183	ASN	2.6
1	C	9	GLN	2.5
1	C	180	GLU	2.5
1	B	8	GLY	2.5
1	A	186	THR	2.5
1	E	186	THR	2.4
1	B	183	ASN	2.4
1	C	11	VAL	2.3
1	C	5	ASN	2.3
1	F	179	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	186	THR	2.2
1	F	183	ASN	2.2
1	B	7	GLN	2.2
1	B	4	GLN	2.1
1	E	187	GLU	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 monosaccharides 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.