



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 10, 2024 – 07:11 pm BST

PDB ID : 8R7Y  
Title : Deoxyribonucleoside regulator DeoR in complex with the DNA operator  
Authors : Pachl, P.; Soltysova, M.; Rezacova, P.  
Deposited on : 2023-11-27  
Resolution : 3.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 2.36.2  
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) : 2.36.2

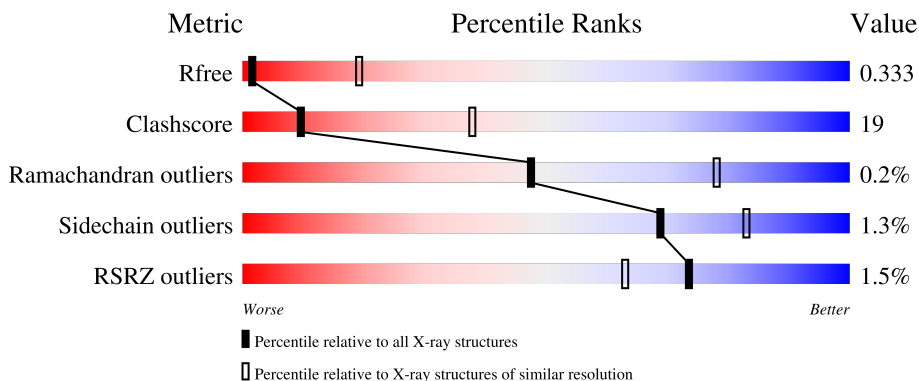
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.70 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.



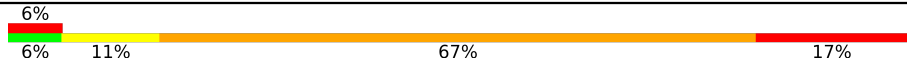
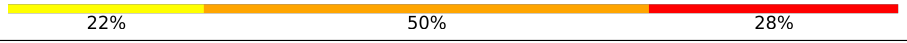
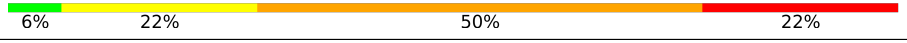
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1049 (3.88-3.52)
Clashscore	141614	1027 (3.86-3.54)
Ramachandran outliers	138981	1069 (3.88-3.52)
Sidechain outliers	138945	1065 (3.88-3.52)
RSRZ outliers	127900	1578 (3.90-3.50)

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	318	 66% 31% 3% 2%
1	B	318	 74% 24% 2% 1%
1	C	318	 63% 34% 3% 2%
1	D	318	 69% 28% 3% 2%
2	E	18	 11% 17% 33% 39%

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Mol	Chain	Length	Quality of chain
2	H	18	 6% 6% 11% 67% 17%
3	F	18	 22% 50% 28%
3	G	18	 6% 22% 50% 22%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 11264 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 Deoxyribonucleoside regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	312	2454	1546	429	470	9	0	0	0
1	B	311	2450	1547	428	466	9	0	0	0
1	C	312	2435	1539	419	468	9	0	0	0
1	D	312	2453	1547	428	470	8	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP P39140
A	-3	ILE	-	expression tag	UNP P39140
A	-2	ASP	-	expression tag	UNP P39140
A	-1	PRO	-	expression tag	UNP P39140
A	0	PHE	-	expression tag	UNP P39140
A	1	THR	-	expression tag	UNP P39140
B	-4	GLY	-	expression tag	UNP P39140
B	-3	ILE	-	expression tag	UNP P39140
B	-2	ASP	-	expression tag	UNP P39140
B	-1	PRO	-	expression tag	UNP P39140
B	0	PHE	-	expression tag	UNP P39140
B	1	THR	-	expression tag	UNP P39140
C	-4	GLY	-	expression tag	UNP P39140
C	-3	ILE	-	expression tag	UNP P39140
C	-2	ASP	-	expression tag	UNP P39140
C	-1	PRO	-	expression tag	UNP P39140
C	0	PHE	-	expression tag	UNP P39140
C	1	THR	-	expression tag	UNP P39140
D	-4	GLY	-	expression tag	UNP P39140
D	-3	ILE	-	expression tag	UNP P39140
D	-2	ASP	-	expression tag	UNP P39140

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	PRO	-	expression tag	UNP P39140
D	0	PHE	-	expression tag	UNP P39140
D	1	THR	-	expression tag	UNP P39140

- Molecule 2 is a DNA chain called OL18 DNA operator, strand 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	E	18	Total	C	N	O	P	0	0	0
			366	179	61	109	17			
2	H	18	Total	C	N	O	P	0	0	0
			366	179	61	109	17			

- Molecule 3 is a DNA chain called OL18 DNA operator, strand 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	F	18	Total	C	N	O	P	0	0	0
			366	178	68	103	17			
3	G	18	Total	C	N	O	P	0	0	0
			366	178	68	103	17			

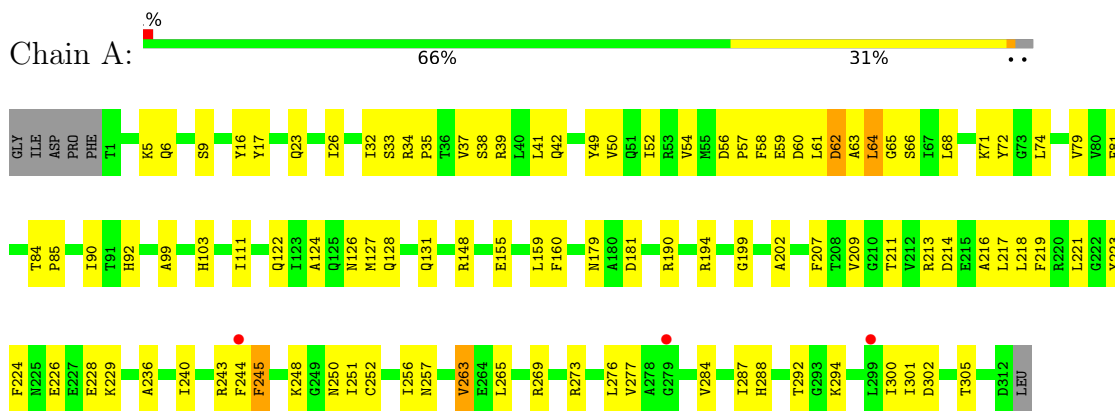
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		
4	B	1	Total	O	0	0
			1	1		
4	C	1	Total	O	0	0
			1	1		
4	E	1	Total	O	0	0
			1	1		
4	G	1	Total	O	0	0
			1	1		
4	H	2	Total	O	0	0
			2	2		

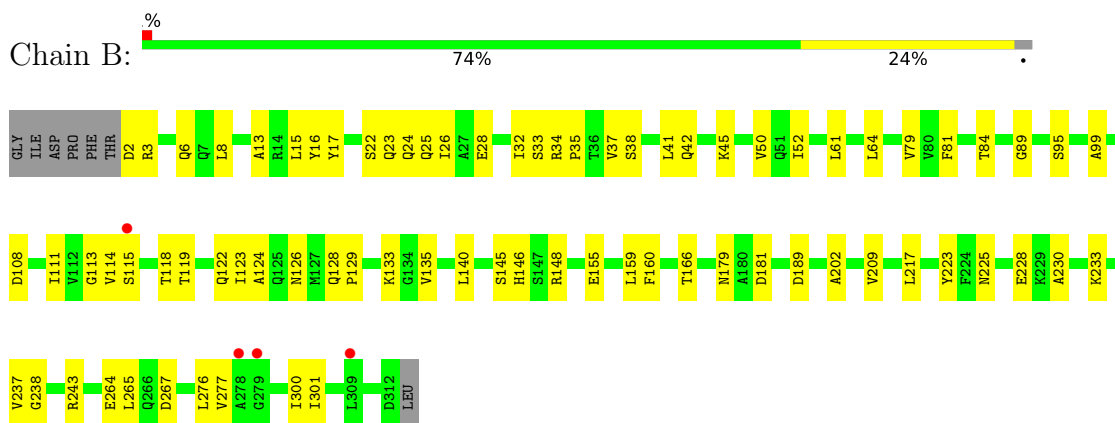
### 3 Residue-property plots [i](#)

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.

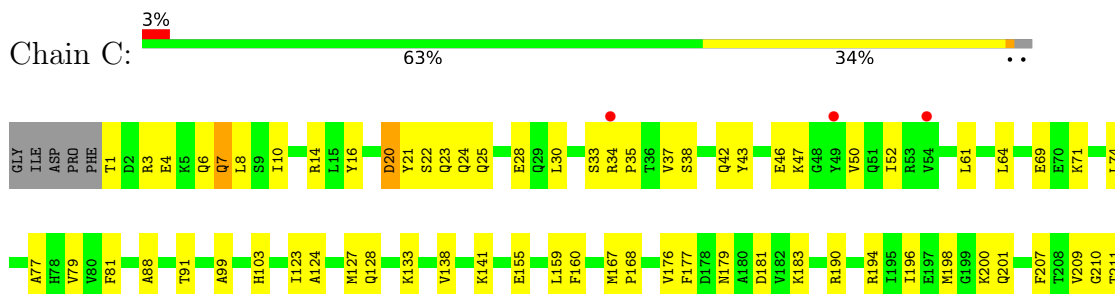
- Molecule 1: Deoxyribonucleoside regulator

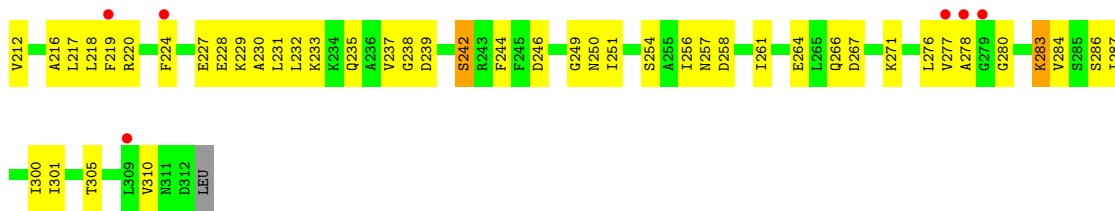


- Molecule 1: Deoxyribonucleoside regulator

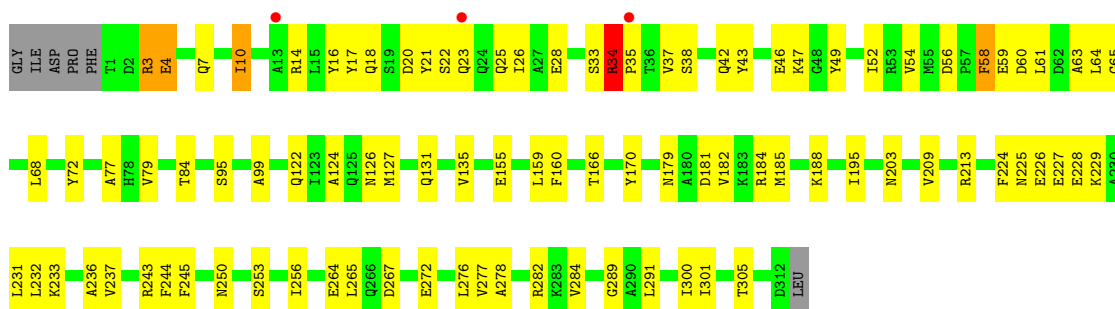


- Molecule 1: Deoxyribonucleoside regulator





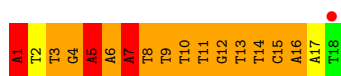
- Molecule 1: Deoxyribonucleoside regulator



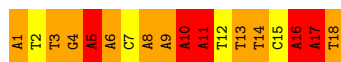
- Molecule 2: OL18 DNA operator, strand 1



- Molecule 2: OL18 DNA operator, strand 1



- Molecule 3: OL18 DNA operator, strand 2



- Molecule 3: OL18 DNA operator, strand 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	166.67Å 166.67Å 331.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.92 – 3.70 48.92 – 3.70	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.92-3.70) 99.3 (48.92-3.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.01 (at 3.67Å)	Xtrriage
Refinement program	REFMAC 5.8.0405	Depositor
R, $R_{free}$	0.266 , 0.333 0.266 , 0.333	Depositor DCC
$R_{free}$ test set	1482 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	149.4	Xtrriage
Anisotropy	0.055	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 109.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.38$ , $\langle L^2 \rangle = 0.20$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	11264	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	173.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% 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	0.26	0/2493	0.57	0/3367
1	B	0.26	0/2490	0.59	0/3362
1	C	0.26	0/2475	0.58	0/3347
1	D	0.26	0/2493	0.61	0/3369
2	E	0.73	0/409	2.02	21/630 (3.3%)
2	H	0.71	0/409	2.08	19/630 (3.0%)
3	F	0.73	0/411	2.21	25/632 (4.0%)
3	G	0.74	0/411	2.10	21/632 (3.3%)
All	All	0.37	0/11591	0.99	86/15969 (0.5%)

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	D	0	1
2	E	0	8
2	H	0	7
3	F	0	10
3	G	0	10
All	All	0	36

There are no bond length outliers.

The worst 5 of 86 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1	DA	N9-C4-C5	-11.15	101.34	105.80
3	F	18	DT	O5'-P-OP2	-9.70	96.97	105.70
3	G	9	DA	N9-C4-C5	-9.39	102.04	105.80
3	F	17	DA	N9-C4-C5	-9.15	102.14	105.80
3	F	18	DT	N3-C2-O2	-9.05	116.87	122.30

There are no chirality outliers.

5 of 36 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	34	ARG	Sidechain
2	E	1	DA	Sidechain
2	E	4	DG	Sidechain
2	E	5	DA	Sidechain
2	E	6	DA	Sidechain

## 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	2454	0	2456	103	1
1	B	2450	0	2451	70	0
1	C	2435	0	2417	117	0
1	D	2453	0	2445	98	0
2	E	366	0	209	20	2
2	H	366	0	209	20	1
3	F	366	0	206	19	0
3	G	366	0	206	10	1
4	A	2	0	0	0	0
4	B	1	0	0	1	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
4	G	1	0	0	0	0
4	H	2	0	0	0	0
All	All	11264	0	10599	424	3

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 19.

The worst 5 of 424 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:C:23:GLN:OE1	1:C:38:SER:OG	1.69	1.08
1:B:111:ILE:HG22	1:B:202:ALA:HA	1.37	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:52:ILE:CG2	1:D:52:ILE:HD12	1.90	1.02
1:B:23:GLN:OE1	1:B:38:SER:OG	1.76	1.01
1:C:8:LEU:HD11	1:C:30:LEU:HD22	1.42	1.01

All (3) 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)
1:A:213:ARG:NH2	2:E:16:DA:C5'[12_554]	1.54	0.66
3:G:1:DA:O5'	2:H:17:DA:OP2[11_454]	1.60	0.60
2:E:18:DT:O2	2:E:18:DT:O2[10_445]	2.15	0.05

## 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	310/318 (98%)	300 (97%)	8 (3%)	2 (1%)	25	62
1	B	309/318 (97%)	300 (97%)	9 (3%)	0	100	100
1	C	310/318 (98%)	299 (96%)	11 (4%)	0	100	100
1	D	310/318 (98%)	299 (96%)	10 (3%)	1 (0%)	41	74
All	All	1239/1272 (97%)	1198 (97%)	38 (3%)	3 (0%)	47	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	ASP
1	A	64	LEU
1	D	4	GLU

### 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	263/271 (97%)	261 (99%)	2 (1%)	81	89
1	B	262/271 (97%)	261 (100%)	1 (0%)	91	95
1	C	259/271 (96%)	254 (98%)	5 (2%)	57	76
1	D	262/271 (97%)	256 (98%)	6 (2%)	50	71
All	All	1046/1084 (96%)	1032 (99%)	14 (1%)	69	83

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	283	LYS
1	D	3	ARG
1	D	84	THR
1	D	34	ARG
1	D	58	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	250	ASN
1	A	257	ASN
1	C	257	ASN
1	D	158	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	312/318 (98%)	-0.21	3 (0%) 82 73	107, 176, 231, 265	0
1	B	311/318 (97%)	-0.18	4 (1%) 77 67	96, 165, 222, 250	0
1	C	312/318 (98%)	-0.06	9 (2%) 51 39	98, 194, 242, 275	0
1	D	312/318 (98%)	-0.33	3 (0%) 82 73	96, 160, 209, 268	0
2	E	18/18 (100%)	-0.80	0 100 100	131, 158, 183, 192	0
2	H	18/18 (100%)	-0.40	1 (5%) 24 17	138, 160, 243, 243	0
3	F	18/18 (100%)	-0.71	0 100 100	124, 150, 206, 207	0
3	G	18/18 (100%)	-0.71	0 100 100	111, 150, 192, 206	0
All	All	1319/1344 (98%)	-0.22	20 (1%) 73 63	96, 173, 232, 275	0

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	278	ALA	10.3
1	C	279	GLY	8.2
1	C	224	PHE	2.8
1	C	309	LEU	2.8
1	B	115	SER	2.7

### 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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.