



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 15, 2017 – 08:59 AM EDT

PDB ID : 5BQL  
Title : Fluorescent protein cyOFP  
Authors : Sens, A.; Ataie, N.; Ng, H.L.; Lin, M.Z.; Chu, J.  
Deposited on : unknown  
Resolution : 2.39 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20029824  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20029824

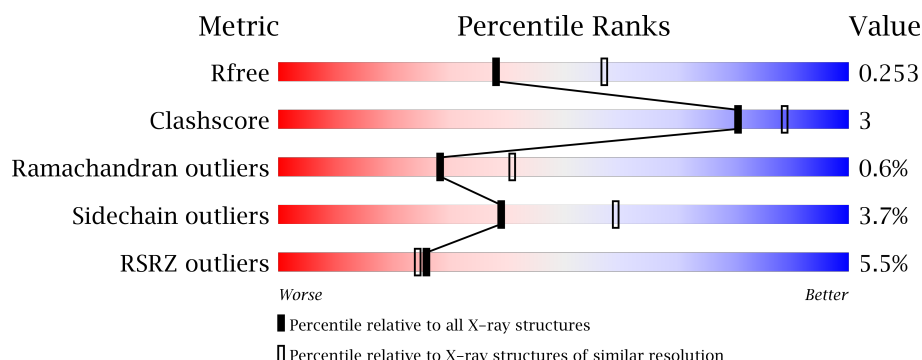
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.39 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))
$R_{free}$	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	242	<div> <div>8%</div> <div> <div></div> <div>84%</div> <div>9%</div> <div>7%</div> </div> </div>
1	B	242	<div> <div>7%</div> <div> <div></div> <div>85%</div> <div>7%</div> <div>7%</div> </div> </div>
1	C	242	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>12%</div> <div>8%</div> </div> </div>
1	D	242	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>9%</div> <div>9%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7216 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 Fluorescent protein cyOFP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	1	0
			1808	1152	308	337	11			
1	B	224	Total	C	N	O	S	0	1	0
			1795	1145	306	333	11			
1	C	223	Total	C	N	O	S	0	3	0
			1805	1150	311	333	11			
1	D	221	Total	C	N	O	S	0	0	0
			1764	1125	301	327	11			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	O	0	0
			5	5		
2	B	5	Total	O	0	0
			5	5		
2	C	17	Total	O	0	0
			17	17		
2	D	17	Total	O	0	0
			17	17		



- Molecule 1: Fluorescent protein cyOFP



LEU  
ASN  
GLY  
MET  
ASP  
GLU  
LEU  
TYR  
LYS

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.12Å 102.76Å 123.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.39 47.39 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.4 (50.00-2.39) 99.4 (47.39-2.39)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.50 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0123	Depositor
R, $R_{free}$	0.193 , 0.252 0.196 , 0.253	Depositor DCC
$R_{free}$ test set	1858 reflections (5.18%)	DCC
Wilson B-factor (Å <sup>2</sup> )	53.7	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 36.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7216	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NRQ

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	0.58	0/1831	0.80	1/2470 (0.0%)
1	B	0.63	0/1818	0.78	0/2453
1	C	0.70	0/1832	0.87	2/2471 (0.1%)
1	D	0.69	0/1784	0.88	0/2408
All	All	0.65	0/7265	0.83	3/9802 (0.0%)

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	2
1	D	0	1
All	All	0	3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	C	161[A]	ARG	NE-CZ-NH2	-5.77	117.42	120.30
1	C	161[B]	ARG	NE-CZ-NH2	-5.77	117.42	120.30
1	A	161	ARG	NE-CZ-NH2	-5.36	117.62	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	189	THR	Peptide

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Mol	Chain	Res	Type	Group
1	B	90	GLU	Peptide
1	D	133	ASN	Peptide

## 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	1808	0	1770	13	0
1	B	1795	0	1761	5	0
1	C	1805	0	1769	17	0
1	D	1764	0	1726	9	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	17	0	0	0	0
2	D	17	0	0	0	0
All	All	7216	0	7026	44	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 (44) 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:A:43:GLN:HE22	1:A:70:SER:HB3	1.48	0.78
1:C:94:TRP:CE2	1:C:110:GLN:HG2	2.22	0.74
1:A:60:ASP:HB3	1:A:166:LEU:HD11	1.72	0.71
1:B:43:GLN:HE22	1:B:70:SER:HB3	1.58	0.69
1:D:43:GLN:HE22	1:D:70:SER:HB3	1.61	0.66
1:B:95:GLU:HG2	1:B:109:THR:HG22	1.82	0.62
1:D:64:THR:HG21	1:D:178:PHE:HE2	1.66	0.61
1:C:80:LEU:HD13	1:C:225:TYR:CZ	2.38	0.59
1:C:94:TRP:NE1	1:C:110:GLN:HG2	2.19	0.57
1:B:80:LEU:HD13	1:B:225:TYR:CZ	2.41	0.55
1:C:219:GLU:OE2	1:C:221:ALA:N	2.41	0.54
1:A:43:GLN:HE22	1:A:70:SER:CB	2.17	0.53
1:D:64:THR:HG22	1:D:96:ARG:NH1	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:60:ASP:HB3	1:D:166:LEU:HD21	1.93	0.51
1:C:39:TYR:O	1:C:75:LYS:HG3	2.12	0.50
1:C:43:GLN:HE22	1:C:70:SER:HB3	1.78	0.48
1:C:82:ASP:OD2	1:C:85:LYS:HD2	2.14	0.48
1:C:162:CYS:HB3	1:C:178:PHE:HB2	1.95	0.47
1:C:151:LEU:HD23	1:C:160:GLY:HA2	1.97	0.47
1:A:125:LEU:C	1:A:125:LEU:HD12	2.35	0.46
1:A:32:HIS:ND1	1:A:32:HIS:C	2.69	0.46
1:A:206:ILE:HD11	1:A:216:GLU:OE2	2.16	0.45
1:D:97:VAL:HG12	1:D:107:THR:HG22	1.99	0.45
1:A:34:GLY:HA3	1:A:45:ASN:HB3	1.99	0.44
1:A:34:GLY:CA	1:A:45:ASN:HB3	2.48	0.44
1:D:203:LEU:HD23	1:D:204:GLU:N	2.32	0.44
1:B:162:CYS:HB3	1:B:178:PHE:HB2	1.99	0.44
1:B:32:HIS:ND1	1:B:32:HIS:C	2.70	0.44
1:C:88:PHE:O	1:C:186:LYS:NZ	2.51	0.44
1:D:203:LEU:HD23	1:D:203:LEU:C	2.38	0.44
1:A:106:LEU:HD21	1:A:125:LEU:HD13	1.98	0.43
1:C:94:TRP:CE2	1:C:110:GLN:CG	2.98	0.43
1:A:162:CYS:HB3	1:A:178:PHE:HB2	2.01	0.43
1:C:55:LEU:HD13	1:C:57:PHE:CZ	2.52	0.43
1:C:86:GLN:HB2	1:C:191:LEU:HD23	2.01	0.43
1:D:64:THR:HG23	1:D:67:NRQ:CB2	2.48	0.43
1:D:32:HIS:C	1:D:32:HIS:ND1	2.73	0.42
1:C:32:HIS:C	1:C:32:HIS:ND1	2.73	0.41
1:C:86:GLN:CB	1:C:191:LEU:HD23	2.50	0.41
1:C:94:TRP:NE1	1:C:110:GLN:CG	2.83	0.41
1:C:88:PHE:HB3	1:C:89:PRO:HA	2.03	0.41
1:A:60:ASP:CB	1:A:166:LEU:HD11	2.46	0.41
1:A:88:PHE:HB3	1:A:89:PRO:HA	2.02	0.40
1:A:67:NRQ:N2	1:A:67:NRQ:HD2	2.36	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	222/242 (92%)	211 (95%)	10 (4%)	1 (0%)	32	46
1	B	220/242 (91%)	215 (98%)	4 (2%)	1 (0%)	32	46
1	C	221/242 (91%)	217 (98%)	2 (1%)	2 (1%)	20	29
1	D	216/242 (89%)	209 (97%)	6 (3%)	1 (0%)	32	46
All	All	879/968 (91%)	852 (97%)	22 (2%)	5 (1%)	28	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	155	ASP
1	B	91	GLY
1	C	24	ASN
1	A	8	LEU
1	C	60	ASP

### 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	193/206 (94%)	186 (96%)	7 (4%)	40	60
1	B	192/206 (93%)	184 (96%)	8 (4%)	34	53
1	C	193/206 (94%)	188 (97%)	5 (3%)	51	72
1	D	188/206 (91%)	180 (96%)	8 (4%)	33	52
All	All	766/824 (93%)	738 (96%)	28 (4%)	39	59

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LEU
1	A	22	SER
1	A	32	HIS

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Mol	Chain	Res	Type
1	A	106	LEU
1	A	122	ASN
1	A	190	ASN
1	A	208	GLU
1	B	11	GLU
1	B	22	SER
1	B	32	HIS
1	B	80	LEU
1	B	163	ASP
1	B	203	LEU
1	B	227	ASP
1	B	230	SER
1	C	32	HIS
1	C	35	GLU
1	C	110	GLN
1	C	216	GLU
1	C	227	ASP
1	D	32	HIS
1	D	44	THR
1	D	116	ASP
1	D	122	ASN
1	D	126	ARG
1	D	180	THR
1	D	189	THR
1	D	226	CYS

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

Mol	Chain	Res	Type
1	A	43	GLN
1	B	26	HIS
1	B	43	GLN
1	B	45	ASN
1	B	190	ASN
1	C	43	GLN
1	C	110	GLN
1	D	32	HIS
1	D	43	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
1	NRQ	A	67	1	23,24,25	3.11	6 (26%)	25,32,34	4.08	8 (32%)
1	NRQ	B	67	1	23,24,25	3.57	4 (17%)	25,32,34	5.46	7 (28%)
1	NRQ	C	67	1	23,24,25	3.13	5 (21%)	25,32,34	5.54	9 (36%)
1	NRQ	D	67	1	23,24,25	3.31	6 (26%)	25,32,34	3.95	5 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	NRQ	A	67	1	-	0/9/31/32	0/2/2/2
1	NRQ	B	67	1	-	0/9/31/32	0/2/2/2
1	NRQ	C	67	1	-	0/9/31/32	0/2/2/2
1	NRQ	D	67	1	-	0/9/31/32	0/2/2/2

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	67	NRQ	CA2-C2	-5.82	1.42	1.48
1	C	67	NRQ	CA2-C2	-4.81	1.43	1.48
1	A	67	NRQ	CA2-C2	-3.83	1.44	1.48
1	C	67	NRQ	C2-N3	-3.69	1.30	1.39
1	B	67	NRQ	C2-N3	-3.67	1.30	1.39
1	D	67	NRQ	CA2-C2	-3.63	1.44	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	67	NRQ	C2-N3	-3.43	1.31	1.39
1	D	67	NRQ	C2-N3	-3.30	1.31	1.39
1	C	67	NRQ	CA2-N2	-2.89	1.32	1.38
1	A	67	NRQ	CA2-N2	-2.52	1.32	1.38
1	D	67	NRQ	CA2-N2	-2.43	1.33	1.38
1	A	67	NRQ	C1-N2	2.08	1.37	1.33
1	B	67	NRQ	C1-N2	2.51	1.38	1.33
1	A	67	NRQ	O2-C2	2.63	1.28	1.23
1	C	67	NRQ	C1-N2	2.86	1.38	1.33
1	D	67	NRQ	C1-N2	3.39	1.39	1.33
1	D	67	NRQ	O2-C2	3.88	1.31	1.23
1	C	67	NRQ	CB2-CA2	12.60	1.46	1.35
1	A	67	NRQ	CB2-CA2	13.02	1.46	1.35
1	D	67	NRQ	CB2-CA2	13.81	1.47	1.35
1	B	67	NRQ	CB2-CA2	15.07	1.48	1.35

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	67	NRQ	O2-C2-CA2	-19.62	120.39	130.97
1	B	67	NRQ	O2-C2-CA2	-17.53	121.52	130.97
1	D	67	NRQ	O2-C2-CA2	-11.94	124.53	130.97
1	A	67	NRQ	O2-C2-CA2	-11.04	125.02	130.97
1	C	67	NRQ	O3-C3-CA3	-5.25	109.04	126.38
1	B	67	NRQ	CG2-CB2-CA2	-4.70	124.73	130.19
1	A	67	NRQ	O3-C3-CA3	-4.47	111.60	126.38
1	B	67	NRQ	O3-C3-CA3	-4.37	111.93	126.38
1	B	67	NRQ	C2-CA2-N2	-4.14	105.90	108.93
1	C	67	NRQ	CG2-CB2-CA2	-4.02	125.52	130.19
1	D	67	NRQ	O3-C3-CA3	-3.89	113.53	126.38
1	C	67	NRQ	C2-CA2-N2	-3.87	106.10	108.93
1	A	67	NRQ	C2-CA2-N2	-3.66	106.25	108.93
1	A	67	NRQ	CG2-CB2-CA2	-3.29	126.37	130.19
1	C	67	NRQ	CB1-CA1-N1	-3.09	114.67	125.34
1	D	67	NRQ	CB1-CA1-N1	-2.56	116.51	125.34
1	A	67	NRQ	CB1-CA1-N1	-2.02	118.38	125.34
1	C	67	NRQ	O2-C2-N3	2.01	128.75	124.49
1	B	67	NRQ	CE-SD-CG1	2.05	107.72	100.35
1	C	67	NRQ	CD1-CG2-CD2	2.06	120.72	117.63
1	A	67	NRQ	CA3-N3-C1	2.20	132.62	128.19
1	C	67	NRQ	CE-SD-CG1	2.31	108.64	100.35
1	D	67	NRQ	CA3-N3-C1	2.64	133.50	128.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	67	NRQ	CA3-N3-C1	3.11	134.44	128.19
1	A	67	NRQ	CB2-CA2-C2	3.55	126.79	122.32
1	D	67	NRQ	CA2-C2-N3	13.96	109.53	103.30
1	A	67	NRQ	CA2-C2-N3	14.42	109.73	103.30
1	C	67	NRQ	CA2-C2-N3	16.90	110.84	103.30
1	B	67	NRQ	CA2-C2-N3	18.34	111.48	103.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	67	NRQ	1	0
1	D	67	NRQ	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates 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	225/242 (92%)	0.61	20 (8%) 10 9	48, 73, 104, 125	0
1	B	223/242 (92%)	0.48	17 (7%) 15 13	42, 69, 103, 126	0
1	C	222/242 (91%)	0.11	8 (3%) 43 42	37, 51, 69, 81	0
1	D	220/242 (90%)	0.11	4 (1%) 69 66	39, 54, 75, 87	0
All	All	890/968 (91%)	0.33	49 (5%) 26 24	37, 60, 97, 126	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	7	GLU	5.5
1	A	50	VAL	4.9
1	A	118	CYS	4.5
1	A	10	LYS	4.4
1	A	142	LEU	4.2
1	D	188	ALA	4.0
1	B	191	LEU	3.9
1	B	190	ASN	3.7
1	B	189	THR	3.6
1	B	188	ALA	3.5
1	A	189	THR	3.3
1	C	105	VAL	3.2
1	B	162	CYS	3.0
1	B	117	GLY	3.0
1	C	99	VAL	2.9
1	B	176	CYS	2.8
1	D	162	CYS	2.8
1	A	162	CYS	2.8
1	B	114	LEU	2.7
1	A	54	PRO	2.7
1	B	87	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	97	VAL	2.7
1	C	176	CYS	2.7
1	A	191	LEU	2.7
1	A	155	ASP	2.6
1	D	99	VAL	2.6
1	A	18	TYR	2.6
1	C	162	CYS	2.5
1	A	48	LYS	2.5
1	B	99	VAL	2.5
1	A	114	LEU	2.5
1	D	116	ASP	2.5
1	B	11	GLU	2.4
1	A	117	GLY	2.4
1	B	50	VAL	2.4
1	C	18	TYR	2.3
1	C	161[A]	ARG	2.3
1	A	116	ASP	2.3
1	A	16	LYS	2.3
1	A	115	GLN	2.3
1	B	90	GLU	2.2
1	C	14[A]	HIS	2.1
1	A	79	GLY	2.1
1	B	192	LYS	2.1
1	A	178	PHE	2.0
1	B	178	PHE	2.0
1	C	104	GLY	2.0
1	A	106	LEU	2.0
1	B	13	MET	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	NRQ	B	67	23/24	0.96	0.12	-	50,54,69,76	0
1	NRQ	C	67	23/24	0.97	0.11	-	37,43,50,55	0
1	NRQ	A	67	23/24	0.95	0.13	-	57,63,75,78	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
1	NRQ	D	67	23/24	0.96	0.12	-	40,42,54,58	0

### 6.3 Carbohydrates [i](#)

There are no carbohydrates 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.