



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 5, 2023 – 06:21 PM EST

PDB ID : 7SFA  
Title : Branchiostoma floridae fluorescent protein LanFP10A2  
Authors : Pletnev, S.; Pletneva, N.; Pletnev, V.Z.; Muslinkina, L.  
Deposited on : 2021-10-03  
Resolution : 1.65 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  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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

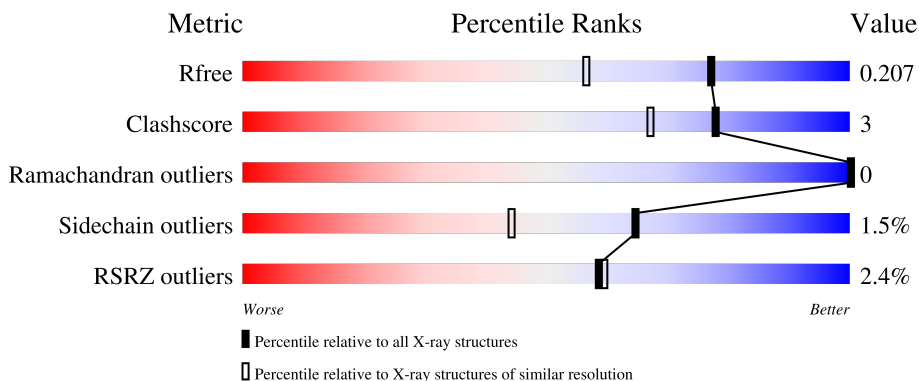
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.65 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	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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	236	 85% 6% 8%
1	B	236	 87% 6% 6%
1	C	236	 85% 7% 8%
1	D	236	 84% 7% 8%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7834 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 lanFP10A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	216	Total 1785	C 1140	N 297	O 337	S 11	0	6	0
1	B	222	Total 1825	C 1163	N 308	O 343	S 11	0	6	0
1	C	217	Total 1772	C 1132	N 295	O 334	S 11	0	4	0
1	D	218	Total 1792	C 1142	N 299	O 340	S 11	0	6	0

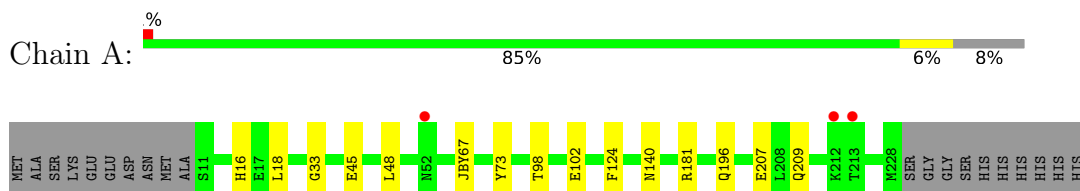
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	163	Total 163	O 163	0	0
2	B	178	Total 178	O 178	0	0
2	C	175	Total 175	O 175	0	0
2	D	144	Total 144	O 144	0	0

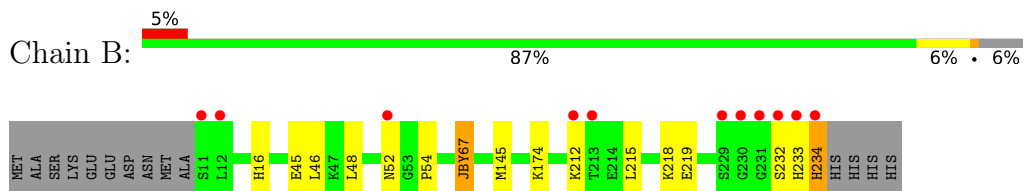
### 3 Residue-property plots [i](#)

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.

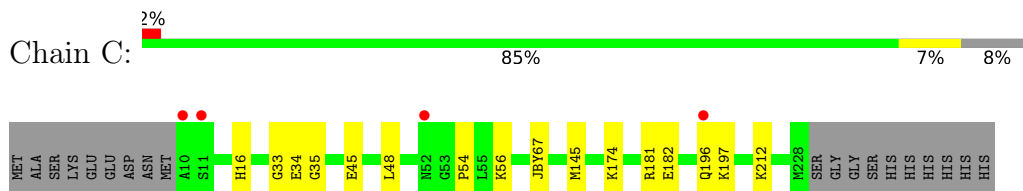
- Molecule 1: Fluorescent protein lanFP10A2



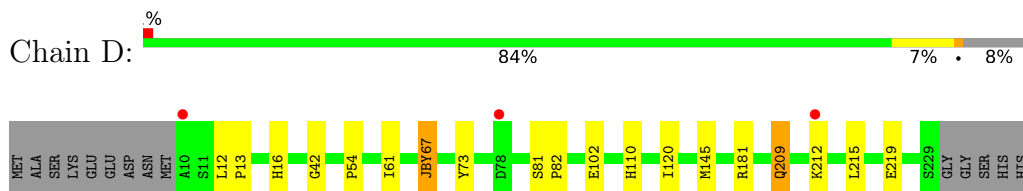
- Molecule 1: Fluorescent protein lanFP10A2



- Molecule 1: Fluorescent protein lanFP10A2



- Molecule 1: Fluorescent protein lanFP10A2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.24Å 96.22Å 158.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.79 – 1.65 27.77 – 1.65	Depositor EDS
% Data completeness (in resolution range)	97.9 (27.79-1.65) 97.9 (27.77-1.65)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.15 (at 1.65Å)	Xtrriage
Refinement program	REFMAC 5.8.0222	Depositor
R, $R_{free}$	0.165 , 0.197 0.177 , 0.207	Depositor DCC
$R_{free}$ test set	1096 reflections (1.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	15.4	Xtrriage
Anisotropy	0.194	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 47.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7834	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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: JBY

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	0.77	1/1814 (0.1%)	0.86	0/2449
1	B	0.73	0/1856	0.87	0/2504
1	C	0.75	2/1801 (0.1%)	0.87	0/2433
1	D	0.73	1/1821 (0.1%)	0.85	0/2460
All	All	0.75	4/7292 (0.1%)	0.86	0/9846

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	102	GLU	CD-OE1	-8.43	1.16	1.25
1	D	102	GLU	CD-OE1	-7.21	1.17	1.25
1	C	182	GLU	CD-OE1	-5.68	1.19	1.25
1	C	182	GLU	CD-OE2	-5.21	1.20	1.25

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	232	SER	Mainchain

## 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	1785	0	1698	8	0
1	B	1825	0	1735	12	0
1	C	1772	0	1686	6	0
1	D	1792	0	1700	14	0
2	A	163	0	0	2	0
2	B	178	0	0	3	0
2	C	175	0	0	4	0
2	D	144	0	0	2	0
All	All	7834	0	6819	40	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 (40) 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:48:LEU:HD13	1:B:215:LEU:HD11	1.53	0.91
1:B:233:HIS:O	1:B:234:HIS:HB2	1.71	0.87
1:A:98:THR:HG21	2:C:439:HOH:O	1.86	0.75
1:A:33:GLY:HA3	1:A:48:LEU:HD23	1.74	0.69
1:C:181[B]:ARG:NH2	2:C:301:HOH:O	2.27	0.67
1:B:48:LEU:CD1	1:B:215:LEU:HD11	2.26	0.64
1:B:48:LEU:HD13	1:B:215:LEU:CD1	2.30	0.58
1:A:45[A]:GLU:HG3	2:A:338:HOH:O	2.04	0.57
1:A:16:HIS:HE1	2:A:324:HOH:O	1.93	0.52
1:B:215:LEU:HD12	1:B:215:LEU:C	2.31	0.50
1:B:174:LYS:HE2	2:B:377:HOH:O	2.12	0.50
1:D:54:PRO:HB3	1:D:212:LYS:HB3	1.96	0.47
1:C:16:HIS:HE1	2:C:342:HOH:O	1.97	0.47
1:B:54:PRO:HB3	1:B:212:LYS:HB3	1.97	0.46
1:A:18:LEU:HD11	1:A:124:PHE:CD2	2.51	0.46
1:D:16:HIS:HE1	2:D:308:HOH:O	1.98	0.46
1:C:54:PRO:HB3	1:C:212:LYS:HB3	1.97	0.45
1:B:45[B]:GLU:HG2	1:B:218:LYS:HG2	1.99	0.45
1:B:52:ASN:ND2	2:B:305:HOH:O	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:181[A]:ARG:NH2	1:D:181[A]:ARG:HG2	2.32	0.45
1:A:207[B]:GLU:OE2	1:A:209:GLN:NE2	2.50	0.44
1:D:215:LEU:C	1:D:215:LEU:HD12	2.38	0.44
1:A:181[B]:ARG:HG2	1:A:181[B]:ARG:HH21	1.84	0.43
1:D:181[A]:ARG:HG2	1:D:181[A]:ARG:HH21	1.83	0.43
1:B:16:HIS:HE1	2:B:320:HOH:O	2.01	0.43
1:D:42:GLY:HA2	1:D:73:TYR:O	2.18	0.43
1:C:56:LYS:HD3	2:C:473:HOH:O	2.18	0.42
1:B:67:JBY:N2	1:B:219:GLU:OE2	2.53	0.42
1:D:81:SER:HB2	1:D:82:PRO:HD2	2.02	0.42
1:D:110:HIS:HE1	2:D:339:HOH:O	2.03	0.41
1:D:16:HIS:CD2	1:D:73:TYR:OH	2.73	0.41
1:D:209:GLN:HE21	1:D:209:GLN:HB2	1.51	0.41
1:C:35:GLY:HA3	1:C:45:GLU:O	2.21	0.41
1:D:67:JBY:N2	1:D:219:GLU:OE2	2.54	0.41
1:B:46:LEU:HD23	1:B:48:LEU:HD11	2.03	0.41
1:C:33:GLY:HA3	1:C:48:LEU:HD23	2.02	0.41
1:D:61:ILE:HD12	1:D:61:ILE:HA	1.87	0.41
1:A:16:HIS:CD2	1:A:73:TYR:OH	2.73	0.41
1:D:12:LEU:HB3	1:D:13:PRO:HD2	2.04	0.40
1:D:13:PRO:HB3	1:D:120:ILE:HD11	2.02	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	217/236 (92%)	215 (99%)	2 (1%)	0	100	100
1	B	223/236 (94%)	222 (100%)	1 (0%)	0	100	100
1	C	216/236 (92%)	214 (99%)	2 (1%)	0	100	100
1	D	219/236 (93%)	218 (100%)	1 (0%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	875/944 (93%)	869 (99%)	6 (1%)	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	193/203 (95%)	191 (99%)	2 (1%)	76	62
1	B	197/203 (97%)	195 (99%)	2 (1%)	76	62
1	C	191/203 (94%)	186 (97%)	5 (3%)	46	21
1	D	194/203 (96%)	192 (99%)	2 (1%)	76	62
All	All	775/812 (95%)	764 (99%)	11 (1%)	65	46

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	140	ASN
1	A	196	GLN
1	B	145	MET
1	B	234	HIS
1	C	34	GLU
1	C	145	MET
1	C	174	LYS
1	C	196	GLN
1	C	197	LYS
1	D	145	MET
1	D	209	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	HIS

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Mol	Chain	Res	Type
1	A	88	HIS
1	A	110	HIS
1	A	140	ASN
1	B	16	HIS
1	B	52	ASN
1	B	110	HIS
1	B	210	HIS
1	C	16	HIS
1	C	196	GLN
1	C	209	GLN
1	D	16	HIS
1	D	110	HIS
1	D	209	GLN
1	D	210	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](#)

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	JBY	A	67	1	19,21,22	3.73	3 (15%)	24,29,31	2.93	10 (41%)
1	JBY	D	67	1	19,21,22	4.31	3 (15%)	24,29,31	2.99	11 (45%)
1	JBY	B	67	1	19,21,22	3.75	7 (36%)	24,29,31	3.57	10 (41%)
1	JBY	C	67	1	19,21,22	4.09	7 (36%)	24,29,31	3.36	11 (45%)

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	JBY	A	67	1	-	0/8/28/30	0/2/2/2
1	JBY	D	67	1	-	0/8/28/30	0/2/2/2
1	JBY	B	67	1	-	0/8/28/30	0/2/2/2
1	JBY	C	67	1	-	0/8/28/30	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	67	JBY	CB2-CA2	15.98	1.48	1.35
1	A	67	JBY	CB2-CA2	14.74	1.47	1.35
1	C	67	JBY	CB2-CA2	14.57	1.47	1.35
1	B	67	JBY	CB2-CA2	12.81	1.45	1.35
1	D	67	JBY	CA2-C2	-8.45	1.40	1.48
1	C	67	JBY	CA2-C2	-7.45	1.41	1.48
1	B	67	JBY	CA2-C2	-7.27	1.41	1.48
1	C	67	JBY	C1-N2	4.46	1.40	1.32
1	A	67	JBY	O2-C2	4.11	1.31	1.23
1	B	67	JBY	CA1-C1	3.73	1.53	1.49
1	D	67	JBY	OH-CZ	-3.32	1.29	1.37
1	A	67	JBY	OH-CZ	-3.22	1.29	1.37
1	B	67	JBY	O2-C2	2.70	1.28	1.23
1	C	67	JBY	O2-C2	2.53	1.28	1.23
1	B	67	JBY	C1-N2	2.49	1.37	1.32
1	C	67	JBY	CD1-CE1	2.43	1.43	1.38
1	C	67	JBY	OH-CZ	-2.36	1.31	1.37
1	C	67	JBY	CA2-N2	-2.36	1.33	1.38
1	B	67	JBY	OH-CZ	-2.22	1.31	1.37
1	B	67	JBY	C1-N3	2.11	1.41	1.37

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	67	JBY	CA2-C2-N3	12.91	109.34	103.39
1	B	67	JBY	CA2-C2-N3	11.99	108.92	103.39
1	D	67	JBY	CA2-C2-N3	8.42	107.27	103.39
1	A	67	JBY	O2-C2-CA2	-7.94	126.50	130.96
1	A	67	JBY	CA2-C2-N3	7.06	106.65	103.39
1	B	67	JBY	O2-C2-CA2	-5.87	127.66	130.96
1	B	67	JBY	CB2-CA2-C2	-5.06	116.24	122.28
1	D	67	JBY	O2-C2-CA2	-4.93	128.19	130.96
1	D	67	JBY	CB3-CA3-N3	-4.50	104.77	111.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	67	JBY	C1-CA1-N1	-4.47	102.95	112.85
1	D	67	JBY	CB2-CA2-C2	-4.44	116.97	122.28
1	C	67	JBY	N3-C1-N2	-4.35	110.79	113.67
1	B	67	JBY	CB3-CA3-N3	-4.33	105.01	111.05
1	B	67	JBY	CB2-CA2-N2	4.17	134.61	128.83
1	A	67	JBY	C1-CA1-N1	-4.05	103.89	112.85
1	B	67	JBY	C1-CA1-N1	-3.97	104.07	112.85
1	A	67	JBY	CB3-CA3-N3	-3.89	105.62	111.05
1	B	67	JBY	CD1-CG2-CD2	3.60	122.96	117.64
1	D	67	JBY	CB2-CA2-N2	3.52	133.70	128.83
1	C	67	JBY	CB2-CA2-C2	-3.43	118.18	122.28
1	C	67	JBY	O2-C2-CA2	-3.30	129.11	130.96
1	C	67	JBY	C1-CA1-N1	-3.26	105.64	112.85
1	A	67	JBY	CE1-CD1-CG2	-3.15	117.14	121.25
1	C	67	JBY	CB2-CA2-N2	3.14	133.18	128.83
1	B	67	JBY	CE2-CD2-CG2	-3.11	117.20	121.25
1	D	67	JBY	CD1-CG2-CD2	3.07	122.18	117.64
1	D	67	JBY	CE1-CD1-CG2	-3.05	117.27	121.25
1	A	67	JBY	CB2-CA2-N2	2.87	132.81	128.83
1	B	67	JBY	N3-C1-N2	-2.73	111.86	113.67
1	A	67	JBY	N3-C1-N2	-2.57	111.96	113.67
1	D	67	JBY	N3-C1-N2	-2.57	111.97	113.67
1	C	67	JBY	CE2-CD2-CG2	-2.55	117.93	121.25
1	A	67	JBY	CB2-CA2-C2	-2.51	119.27	122.28
1	A	67	JBY	CD1-CG2-CD2	2.50	121.34	117.64
1	C	67	JBY	CD1-CG2-CD2	2.47	121.30	117.64
1	B	67	JBY	C2-N3-C1	-2.38	105.15	108.28
1	A	67	JBY	CD2-CE2-CZ	-2.37	117.27	119.88
1	C	67	JBY	CD2-CE2-CZ	2.36	122.46	119.88
1	C	67	JBY	C2-N3-C1	-2.27	105.29	108.28
1	D	67	JBY	CA1-C1-N3	2.23	126.28	122.89
1	D	67	JBY	CA1-C1-N2	-2.11	121.45	124.28
1	C	67	JBY	CG2-CB2-CA2	-2.06	127.42	129.94

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	D	67	JBY	1	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	67	JBY	1	0

## 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	215/236 (91%)	-0.17	3 (1%) 75 79	9, 15, 31, 69	0
1	B	221/236 (93%)	-0.07	11 (4%) 28 27	9, 16, 32, 55	0
1	C	216/236 (91%)	-0.17	4 (1%) 66 69	9, 16, 32, 48	0
1	D	217/236 (91%)	-0.15	3 (1%) 75 79	10, 18, 33, 46	0
All	All	869/944 (92%)	-0.14	21 (2%) 59 59	9, 16, 33, 69	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	234	HIS	10.0
1	B	230	GLY	7.3
1	B	233	HIS	7.1
1	C	10	ALA	6.3
1	D	10	ALA	4.9
1	B	11	SER	4.9
1	B	212	LYS	4.9
1	A	213	THR	4.7
1	B	232	SER	4.5
1	B	229	SER	4.2
1	B	231	GLY	4.0
1	A	212	LYS	3.9
1	A	52	ASN	3.3
1	C	196	GLN	3.0
1	B	12	LEU	2.4
1	D	78[A]	ASP	2.3
1	C	11	SER	2.3
1	B	213	THR	2.2
1	B	52	ASN	2.2
1	C	52	ASN	2.1
1	D	212	LYS	2.1

## 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
1	JBY	C	67	20/21	0.94	0.14	11,13,20,22	0
1	JBY	D	67	20/21	0.95	0.11	13,15,18,19	0
1	JBY	A	67	20/21	0.96	0.11	11,12,19,20	0
1	JBY	B	67	20/21	0.97	0.09	11,12,18,19	0

## 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.