



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2025 – 05:11 pm BST

PDB ID : 9FER / pdb_00009fer
Title : wasCFP (SG P21) - Directionality of Optical Properties of Fluorescent Proteins
Authors : Myskova, J.; Brynda, J.; Lazar, J.
Deposited on : 2024-05-21
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

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 Phenix2.0rc1
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.43.1

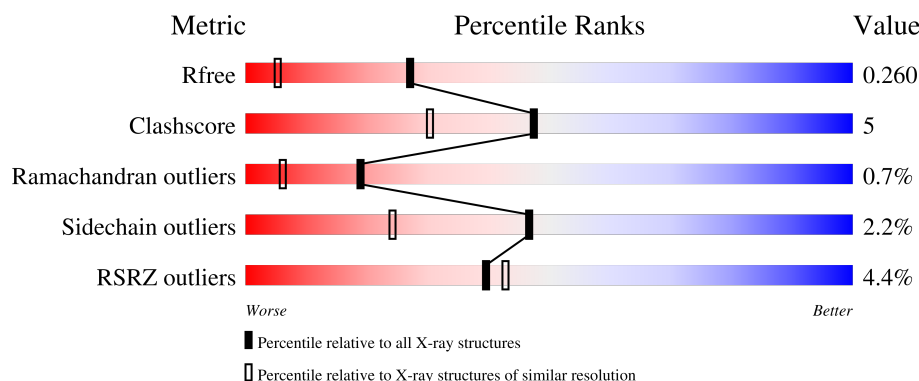
1 Overall quality at a glance

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}	164625	2328 (1.66-1.66)
Clashscore	180529	2515 (1.66-1.66)
Ramachandran outliers	177936	2475 (1.66-1.66)
Sidechain outliers	177891	2475 (1.66-1.66)
RSRZ outliers	164620	2328 (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 ≥ 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	258	<div> <div>4%</div> <div>71%</div> <div>12%</div> <div>•</div> <div>16%</div> </div>
1	B	258	<div> <div>3%</div> <div>73%</div> <div>10%</div> <div>•</div> <div>16%</div> </div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 3816 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 plum.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	2	0
			1745	1107	289	338	11			
1	B	218	Total	C	N	O	S	0	1	0
			1738	1101	289	337	11			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-31	HIS	-	expression tag	UNP Q5S3G7
A	-30	HIS	-	expression tag	UNP Q5S3G7
A	-29	HIS	-	expression tag	UNP Q5S3G7
A	-28	HIS	-	expression tag	UNP Q5S3G7
A	-27	HIS	-	expression tag	UNP Q5S3G7
A	-26	HIS	-	expression tag	UNP Q5S3G7
A	-25	GLY	-	expression tag	UNP Q5S3G7
A	-24	MET	-	expression tag	UNP Q5S3G7
A	-23	ALA	-	expression tag	UNP Q5S3G7
A	-22	SER	-	expression tag	UNP Q5S3G7
A	-21	MET	-	expression tag	UNP Q5S3G7
A	-20	THR	-	expression tag	UNP Q5S3G7
A	-19	GLY	-	expression tag	UNP Q5S3G7
A	-18	GLY	-	expression tag	UNP Q5S3G7
A	-17	GLN	-	expression tag	UNP Q5S3G7
A	-16	GLN	-	expression tag	UNP Q5S3G7
A	-15	MET	-	expression tag	UNP Q5S3G7
A	-14	GLY	-	expression tag	UNP Q5S3G7
A	-13	ARG	-	expression tag	UNP Q5S3G7
A	-12	ASP	-	expression tag	UNP Q5S3G7
A	-11	LEU	-	expression tag	UNP Q5S3G7
A	-10	TYR	-	expression tag	UNP Q5S3G7
A	-9	GLU	-	expression tag	UNP Q5S3G7
A	-8	ASN	-	expression tag	UNP Q5S3G7
A	-7	LEU	-	expression tag	UNP Q5S3G7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	TYR	-	expression tag	UNP Q5S3G7
A	-5	PHE	-	expression tag	UNP Q5S3G7
A	-4	GLN	-	expression tag	UNP Q5S3G7
A	-3	GLY	-	expression tag	UNP Q5S3G7
A	-2	SER	-	expression tag	UNP Q5S3G7
A	-1	SER	-	expression tag	UNP Q5S3G7
A	68	CH6	-	insertion	UNP Q5S3G7
B	-31	HIS	-	expression tag	UNP Q5S3G7
B	-30	HIS	-	expression tag	UNP Q5S3G7
B	-29	HIS	-	expression tag	UNP Q5S3G7
B	-28	HIS	-	expression tag	UNP Q5S3G7
B	-27	HIS	-	expression tag	UNP Q5S3G7
B	-26	HIS	-	expression tag	UNP Q5S3G7
B	-25	GLY	-	expression tag	UNP Q5S3G7
B	-24	MET	-	expression tag	UNP Q5S3G7
B	-23	ALA	-	expression tag	UNP Q5S3G7
B	-22	SER	-	expression tag	UNP Q5S3G7
B	-21	MET	-	expression tag	UNP Q5S3G7
B	-20	THR	-	expression tag	UNP Q5S3G7
B	-19	GLY	-	expression tag	UNP Q5S3G7
B	-18	GLY	-	expression tag	UNP Q5S3G7
B	-17	GLN	-	expression tag	UNP Q5S3G7
B	-16	GLN	-	expression tag	UNP Q5S3G7
B	-15	MET	-	expression tag	UNP Q5S3G7
B	-14	GLY	-	expression tag	UNP Q5S3G7
B	-13	ARG	-	expression tag	UNP Q5S3G7
B	-12	ASP	-	expression tag	UNP Q5S3G7
B	-11	LEU	-	expression tag	UNP Q5S3G7
B	-10	TYR	-	expression tag	UNP Q5S3G7
B	-9	GLU	-	expression tag	UNP Q5S3G7
B	-8	ASN	-	expression tag	UNP Q5S3G7
B	-7	LEU	-	expression tag	UNP Q5S3G7
B	-6	TYR	-	expression tag	UNP Q5S3G7
B	-5	PHE	-	expression tag	UNP Q5S3G7
B	-4	GLN	-	expression tag	UNP Q5S3G7
B	-3	GLY	-	expression tag	UNP Q5S3G7
B	-2	SER	-	expression tag	UNP Q5S3G7
B	-1	SER	-	expression tag	UNP Q5S3G7
B	68	CH6	-	insertion	UNP Q5S3G7

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	163	Total 166	O 166	0	3
2	B	166	Total 167	O 167	0	1

i

- Molecule 1: Fluorescent protein plum



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.32Å 77.75Å 96.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.14 – 1.65 43.14 – 1.65	Depositor EDS
% Data completeness (in resolution range)	98.9 (43.14-1.65) 98.9 (43.14-1.65)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.209 , 0.253 0.223 , 0.260	Depositor DCC
R_{free} test set	2821 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	14.5	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3816	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.78 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.2459e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CH6

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	1.19	5/1769 (0.3%)	1.26	2/2382 (0.1%)
1	B	1.19	3/1759 (0.2%)	1.24	4/2368 (0.2%)
All	All	1.19	8/3528 (0.2%)	1.25	6/4750 (0.1%)

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 (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	17	HIS	CE1-NE2	6.06	1.38	1.32
1	A	94	GLU	CD-OE1	6.00	1.36	1.25
1	A	83	LEU	C-O	5.93	1.31	1.24
1	A	75	HIS	CE1-NE2	5.89	1.38	1.32
1	B	79	ILE	N-CA	5.55	1.50	1.46
1	B	191	GLY	C-O	5.17	1.28	1.23
1	A	20	GLY	C-O	5.13	1.30	1.23
1	A	17	HIS	CE1-NE2	5.10	1.37	1.32

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	ASP	CA-CB-CG	6.99	119.59	112.60
1	B	169	ASP	CA-CB-CG	6.90	119.50	112.60
1	B	14	PHE	CA-CB-CG	6.17	119.97	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	14	PHE	CA-CB-CG	5.85	119.65	113.80
1	B	168	LYS	CB-CA-C	-5.63	102.26	111.50
1	B	141	MET	CA-C-O	-5.16	115.84	121.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	224	GLY	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	1745	0	1664	22	0
1	B	1738	0	1646	15	0
2	A	166	0	0	1	0
2	B	167	0	0	2	0
All	All	3816	0	3310	37	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 5.

All (37) 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:141:MET:HE2	1:A:168:LYS:HA	1.54	0.88
1:A:141:MET:CE	1:A:168:LYS:HA	2.10	0.81
1:A:141:MET:HE2	1:A:168:LYS:CA	2.26	0.65
1:A:168:LYS:O	1:A:169:ASP:CB	2.47	0.62
1:B:149:ARG:NH2	1:B:160:GLU:OE2	2.26	0.61
1:A:168:LYS:O	1:A:169:ASP:HB2	2.00	0.60
1:A:147:SER:OG	1:A:194:LYS:NZ	2.36	0.59
1:B:199:LEU:HD23	1:B:200:ASP:N	2.20	0.57
1:B:199:LEU:HD23	1:B:199:LEU:C	2.29	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:LEU:C	1:A:199:LEU:HD23	2.30	0.57
1:B:216:ARG:NE	2:B:303:HOH:O	2.38	0.56
1:A:194:LYS:HG2	1:A:222:SER:HB3	1.86	0.56
1:B:169:ASP:HB3	2:B:344:HOH:O	2.05	0.55
1:B:12:MET:HE1	1:B:113:LEU:HD11	1.89	0.54
1:B:68:CH6:HD2	1:B:161:MET:HE1	1.91	0.52
1:A:9:LYS:HG2	2:A:391:HOH:O	2.10	0.51
1:B:92:LYS:NZ	1:B:110:ASP:OD2	2.44	0.50
1:A:199:LEU:HD23	1:A:200:ASP:N	2.28	0.48
1:B:109:GLN:C	1:B:109:GLN:HE21	2.21	0.48
1:B:70:LYS:NZ	1:B:148:GLU:OE1	2.46	0.47
1:A:20:GLY:HA3	1:A:124:VAL:O	2.15	0.46
1:A:12:MET:HE1	1:A:113:LEU:HD11	1.97	0.46
1:A:59:ASP:HB3	1:A:165:LEU:HD21	1.97	0.45
1:A:25:HIS:HE1	1:A:51:GLY:O	2.00	0.44
1:A:204:HIS:HD2	1:A:205:ASN:O	2.00	0.44
1:A:99:PHE:CD1	1:A:175:ALA:HB2	2.53	0.43
1:A:21:SER:HA	1:A:25:HIS:O	2.18	0.43
1:B:46:LEU:HG	1:B:65:ILE:HD13	2.01	0.43
1:A:140:THR:HG21	1:A:165:LEU:HD13	2.01	0.42
1:A:109:GLN:HE21	1:A:109:GLN:C	2.28	0.42
1:A:141:MET:HE1	1:A:168:LYS:HA	2.00	0.42
1:B:149:ARG:HA	1:B:194:LYS:HD3	2.02	0.42
1:B:199:LEU:C	1:B:199:LEU:CD2	2.92	0.42
1:B:59:ASP:HB3	1:B:165:LEU:HD21	2.01	0.42
1:A:113:LEU:HD12	1:A:117:GLU:C	2.46	0.41
1:A:153:GLU:HB3	1:A:158:LYS:HZ3	1.85	0.41
1:B:87:PHE:HB3	1:B:88:PRO:HA	2.02	0.41

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	214/258 (83%)	208 (97%)	5 (2%)	1 (0%)	25	11
1	B	214/258 (83%)	207 (97%)	5 (2%)	2 (1%)	14	3
All	All	428/516 (83%)	415 (97%)	10 (2%)	3 (1%)	19	6

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	169	ASP
1	A	169	ASP
1	B	223	THR

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	181/216 (84%)	175 (97%)	6 (3%)	33	10
1	B	178/216 (82%)	176 (99%)	2 (1%)	70	54
All	All	359/432 (83%)	351 (98%)	8 (2%)	47	24

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	GLN
1	A	164	ARG
1	A	169	ASP
1	A	194	LYS
1	A	213	GLN
1	A	222	SER
1	B	109	GLN
1	B	169	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	HIS
1	A	109	GLN
1	A	114	GLN
1	A	204	HIS
1	B	98	ASN
1	B	109	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	CH6	B	68	1	24,24,25	4.67	6 (25%)	28,32,34	3.60	10 (35%)
1	CH6	A	68	1	24,24,25	3.54	4 (16%)	28,32,34	3.53	9 (32%)

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	CH6	B	68	1	-	6/12/31/32	0/2/2/2
1	CH6	A	68	1	-	2/12/31/32	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	68	CH6	CB2-CA2	18.12	1.50	1.35
1	A	68	CH6	CB2-CA2	12.10	1.45	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	68	CH6	CA2-C2	-8.55	1.40	1.48
1	A	68	CH6	C1-N2	8.37	1.44	1.32
1	B	68	CH6	C1-N2	8.14	1.44	1.32
1	A	68	CH6	CA2-C2	-6.66	1.42	1.48
1	B	68	CH6	O2-C2	4.39	1.32	1.23
1	A	68	CH6	O2-C2	4.26	1.32	1.23
1	B	68	CH6	C2-N3	-3.87	1.30	1.39
1	B	68	CH6	CE2-CZ	2.49	1.43	1.38

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	68	CH6	CA2-C2-N3	15.08	110.50	103.37
1	A	68	CH6	CA2-C2-N3	12.64	109.35	103.37
1	A	68	CH6	O2-C2-CA2	-9.50	125.63	130.96
1	B	68	CH6	O2-C2-CA2	-6.56	127.28	130.96
1	A	68	CH6	CA1-C1-N3	-4.84	118.54	124.85
1	B	68	CH6	CA2-N2-C1	-4.83	102.22	105.77
1	B	68	CH6	C2-N3-C1	-3.86	106.01	107.97
1	A	68	CH6	CG2-CB2-CA2	-3.49	125.66	129.94
1	A	68	CH6	CA1-C1-N2	3.38	130.38	123.56
1	A	68	CH6	C2-N3-C1	-3.23	106.33	107.97
1	B	68	CH6	O3-C3-CA3	-3.04	117.21	126.39
1	A	68	CH6	CE1-CD1-CG2	-2.75	117.66	121.25
1	B	68	CH6	CA3-N3-C1	2.69	130.39	127.16
1	A	68	CH6	CD1-CG2-CD2	2.67	121.60	117.64
1	B	68	CH6	CB2-CA2-C2	-2.64	119.13	122.28
1	A	68	CH6	CD2-CG2-CB2	-2.61	112.33	121.22
1	B	68	CH6	CA1-C1-N3	-2.55	121.52	124.85
1	B	68	CH6	CD2-CG2-CB2	-2.17	113.84	121.22
1	B	68	CH6	CA1-C1-N2	2.03	127.66	123.56

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	68	CH6	C1-CA1-CB1-CG1
1	B	68	CH6	C2-CA2-CB2-CG2
1	B	68	CH6	C3-CA3-N3-C2
1	A	68	CH6	C2-CA2-CB2-CG2
1	B	68	CH6	CB1-CG1-SD-CE
1	B	68	CH6	C3-CA3-N3-C1

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Mol	Chain	Res	Type	Atoms
1	A	68	CH6	C1-CA1-CB1-CG1
1	B	68	CH6	N1-CA1-CB1-CG1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	68	CH6	1	0

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 [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, 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	216/258 (83%)	0.12	11 (5%) 34 37	10, 18, 29, 52	2 (0%)
1	B	217/258 (84%)	0.08	8 (3%) 45 48	11, 17, 28, 82	1 (0%)
All	All	433/516 (83%)	0.10	19 (4%) 39 43	10, 17, 29, 82	3 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	225	ALA	10.9
1	B	223	THR	6.5
1	B	224	GLY	6.1
1	A	172	HIS	4.7
1	B	168	LYS	3.9
1	A	169	ASP	3.3
1	A	224	GLY	3.2
1	B	169	ASP	2.9
1	B	6	GLU	2.8
1	B	222	SER	2.7
1	A	168	LYS	2.7
1	A	223	THR	2.6
1	A	7	VAL	2.3
1	A	170	GLY	2.3
1	A	52	GLY	2.1
1	A	141	MET	2.1
1	A	222	SER	2.1
1	B	113	LEU	2.1
1	A	6	GLU	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, 95th 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 ²)	Q<0.9
1	CH6	B	68	23/24	0.96	0.07	12,16,18,19	0
1	CH6	A	68	23/24	0.97	0.07	13,16,19,20	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.