



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 08:33 pm BST

PDB ID : 6JC5
Title : Crystal structure of the blue fluorescent protein with a Leu-Leu-Gly tri-peptide chromophore derived from the purple chromoprotein of *Stichodactyla haddoni*
Authors : Ko, T.P.; Huang, K.F.; Chang, H.Y.
Deposited on : 2019-01-28
Resolution : 2.05 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

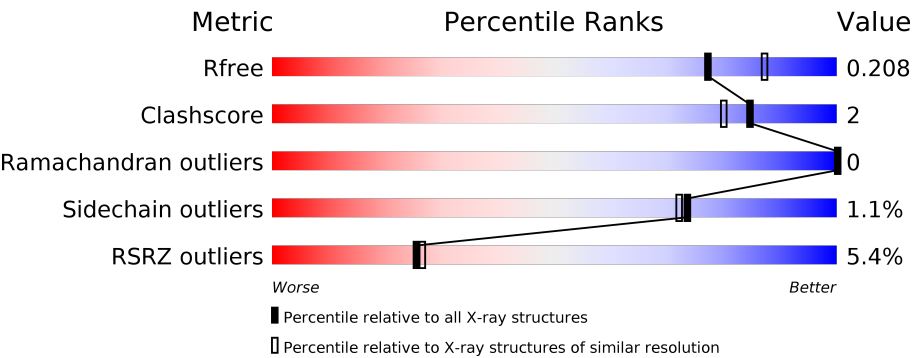
MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

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 2.05 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.




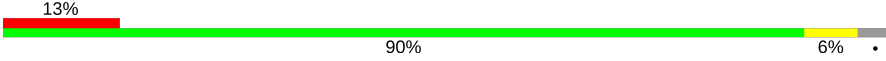
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2684 (2.08-2.04)
Clashscore	141614	2801 (2.08-2.04)
Ramachandran outliers	138981	2768 (2.08-2.04)
Sidechain outliers	138945	2768 (2.08-2.04)
RSRZ outliers	127900	2646 (2.08-2.04)

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 ≥ 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	228	<div><div>2%</div><div><div></div><div>91%</div><div>5%</div><div>.</div></div></div>
1	B	228	<div><div>%</div><div><div></div><div>89%</div><div>7%</div><div>.</div></div></div>
1	C	228	<div><div>3%</div><div><div></div><div>90%</div><div>7%</div><div>.</div></div></div>
1	D	228	<div><div>2%</div><div><div></div><div>92%</div><div>5%</div><div>.</div></div></div>
1	E	228	<div><div>4%</div><div><div></div><div>88%</div><div>8%</div><div>.</div></div></div>
1	F	228	<div><div>4%</div><div><div></div><div>92%</div><div>.</div><div>.</div><div>.</div></div></div>

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Mol	Chain	Length	Quality of chain
1	G	228	 13% 89% 7% •
1	H	228	 13% 90% 6% •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15377 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 shBFP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	3	0
			1777	1134	301	327	15			
1	B	221	Total	C	N	O	S	0	5	0
			1788	1140	304	327	17			
1	C	221	Total	C	N	O	S	0	2	0
			1773	1130	301	326	16			
1	D	221	Total	C	N	O	S	0	6	0
			1794	1144	304	329	17			
1	E	220	Total	C	N	O	S	0	5	0
			1788	1140	304	328	16			
1	F	220	Total	C	N	O	S	0	2	0
			1768	1127	300	326	15			
1	G	221	Total	C	N	O	S	0	0	0
			1761	1121	299	326	15			
1	H	220	Total	C	N	O	S	0	0	0
			1757	1119	298	325	15			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	218	Total	O	0	0
			218	218		
2	B	237	Total	O	0	0
			237	237		
2	C	161	Total	O	0	0
			161	161		
2	D	202	Total	O	0	0
			202	202		
2	E	116	Total	O	0	0
			116	116		
2	F	107	Total	O	0	0
			107	107		

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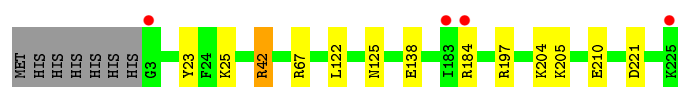
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	74	Total 74	O 74	0	0
2	H	56	Total 56	O 56	0	0

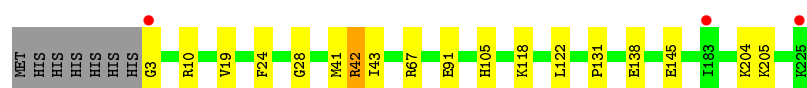
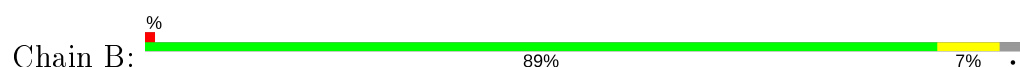
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

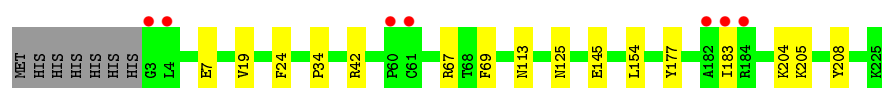
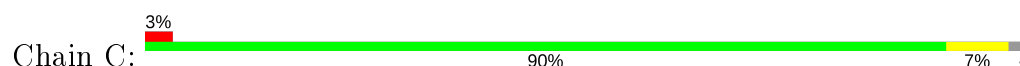
• Molecule 1: shBFP



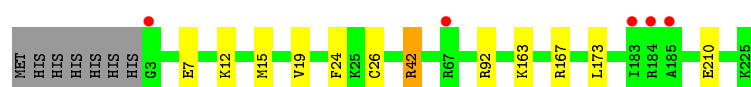
• Molecule 1: shBFP



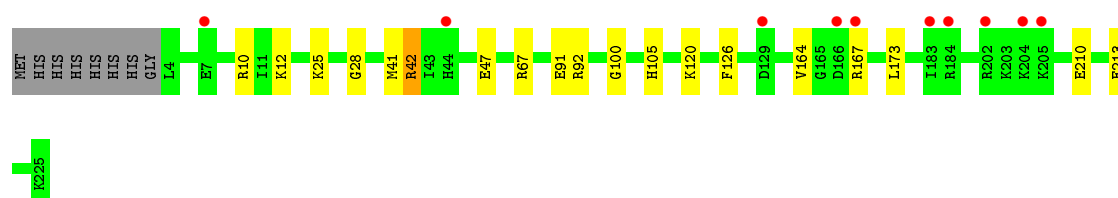
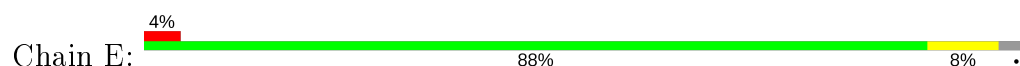
• Molecule 1: shBFP



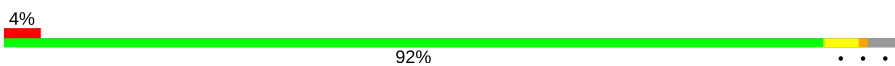
• Molecule 1: shBFP

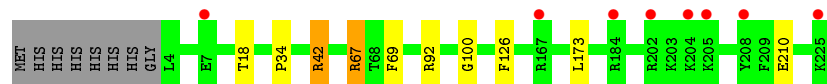


• Molecule 1: shBFP

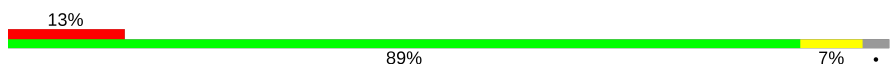


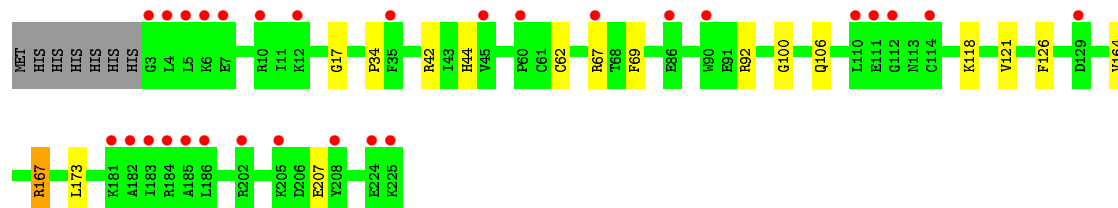
● Molecule 1: shBFP

Chain F: 

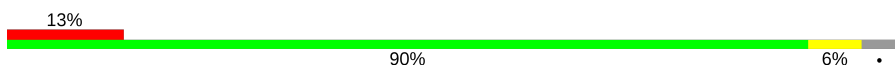


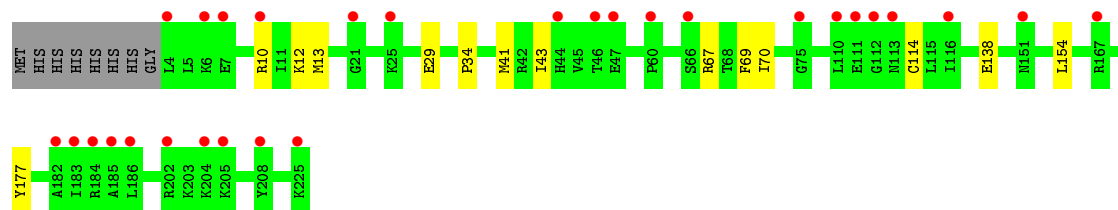
● Molecule 1: shBFP

Chain G: 



● Molecule 1: shBFP

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	103.57Å 98.71Å 197.49Å 90.00° 100.09° 90.00°	Depositor
Resolution (Å)	28.20 – 2.05 29.12 – 2.05	Depositor EDS
% Data completeness (in resolution range)	90.7 (28.20-2.05) 90.7 (29.12-2.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 2.04Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.172 , 0.208 0.172 , 0.208	Depositor DCC
R_{free} test set	2000 reflections (1.72%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 53.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15377	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% of the height of the origin peak. No significant pseudotranslation is detected.*

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: BJJ

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	0.31	0/1812	0.55	0/2443
1	B	0.31	0/1829	0.55	0/2464
1	C	0.30	0/1805	0.54	0/2432
1	D	0.31	0/1838	0.54	0/2476
1	E	0.29	0/1829	0.54	0/2464
1	F	0.28	0/1800	0.52	0/2427
1	G	0.29	0/1786	0.52	0/2407
1	H	0.28	0/1782	0.51	0/2402
All	All	0.30	0/14481	0.53	0/19515

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	G	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	G	62	CYS	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	1777	0	1725	11	0
1	B	1788	0	1741	12	0
1	C	1773	0	1716	8	1
1	D	1794	0	1747	6	1
1	E	1788	0	1738	9	0
1	F	1768	0	1711	5	0
1	G	1761	0	1700	8	0
1	H	1757	0	1697	6	0
2	A	218	0	0	6	0
2	B	237	0	0	4	0
2	C	161	0	0	0	0
2	D	202	0	0	3	0
2	E	116	0	0	0	0
2	F	107	0	0	1	0
2	G	74	0	0	1	0
2	H	56	0	0	0	0
All	All	15377	0	13775	64	1

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 2.

All (64) 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:D:167:ARG:NH2	2:D:301:HOH:O	2.18	0.75
1:B:10[A]:ARG:NH2	2:B:303:HOH:O	2.23	0.70
1:H:67:ARG:HE	1:H:70:ILE:HD11	1.58	0.68
1:F:67:ARG:NH1	2:F:302:HOH:O	2.27	0.67
1:C:183:ILE:HD12	1:C:183:ILE:H	1.61	0.66
1:H:10:ARG:NH2	1:H:29:GLU:OE1	2.30	0.63
1:A:125:ASN:ND2	2:A:304:HOH:O	2.27	0.62
1:B:3:GLY:N	2:B:308:HOH:O	2.33	0.61
1:B:41:MET:SD	1:B:43:ILE:HD11	2.43	0.58
1:E:67[B]:ARG:NH2	1:E:213:GLU:OE2	2.37	0.57
1:A:184:ARG:NH1	2:A:308:HOH:O	2.36	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:LYS:HB3	1:C:208:TYR:HB2	1.85	0.57
1:B:67:ARG:NH2	1:B:145:GLU:OE1	2.33	0.55
1:A:138:GLU:OE2	2:A:301:HOH:O	2.18	0.54
1:B:91:GLU:OE1	2:B:301:HOH:O	2.18	0.54
1:A:42:ARG:HG3	1:A:210:GLU:HG2	1.89	0.54
1:D:42:ARG:HG3	1:D:210:GLU:HG2	1.91	0.53
1:G:17:GLY:HA3	1:G:121:VAL:HG22	1.90	0.53
1:A:221:ASP:OD2	2:A:302:HOH:O	2.19	0.50
1:B:204:LYS:HG2	1:B:205:LYS:HG3	1.92	0.50
1:B:91:GLU:HG2	1:B:105:HIS:HD2	1.76	0.50
1:G:92:ARG:HD3	1:G:173:LEU:HD23	1.94	0.49
1:E:100:GLY:HA3	1:E:126:PHE:CD1	2.47	0.49
1:A:23:TYR:OH	1:A:25:LYS:HE3	2.13	0.49
1:F:42:ARG:HG3	1:F:210:GLU:HG2	1.95	0.49
1:E:42:ARG:HG3	1:E:210:GLU:HG2	1.96	0.48
1:A:138:GLU:OE1	2:A:303:HOH:O	2.20	0.48
1:F:100:GLY:HA3	1:F:126:PHE:CD1	2.49	0.48
1:G:164:VAL:HB	1:G:167:ARG:HG2	1.94	0.47
1:C:67:ARG:NH2	1:C:145:GLU:OE1	2.43	0.47
1:E:92:ARG:HD3	1:E:173:LEU:HD23	1.95	0.47
1:E:91[A]:GLU:HG2	1:E:105:HIS:HD2	1.80	0.47
1:G:67:ARG:NH2	2:G:305:HOH:O	2.37	0.45
1:E:25:LYS:HB2	1:E:47:GLU:HB2	1.98	0.45
1:F:92:ARG:HD3	1:F:173:LEU:HD23	1.98	0.45
1:C:34:PRO:HA	1:C:69:PHE:HA	1.99	0.45
1:C:154:LEU:HB3	1:C:177:TYR:HB2	1.98	0.44
1:G:44:HIS:HE1	1:G:207:GLU:OE2	2.00	0.44
1:H:12:LYS:HG3	1:H:114:CYS:SG	2.58	0.44
1:A:204:LYS:HG3	1:A:205:LYS:HG2	2.00	0.44
1:D:15:MET:HB3	1:D:26[B]:CYS:SG	2.58	0.44
1:H:34:PRO:HA	1:H:69:PHE:HA	2.00	0.44
1:C:125:ASN:HB2	2:D:429:HOH:O	2.18	0.43
1:E:28:GLY:HA3	1:E:41:MET:HE2	2.00	0.43
1:D:92:ARG:HD3	1:D:173:LEU:HD23	2.00	0.43
1:G:34:PRO:HA	1:G:69:PHE:HA	2.01	0.43
1:H:41:MET:SD	1:H:43:ILE:HD11	2.59	0.42
1:D:19:VAL:HG23	1:D:24:PHE:HE1	1.83	0.42
1:E:164:VAL:O	1:E:167:ARG:HG2	2.19	0.42
1:C:19:VAL:HG23	1:C:24:PHE:HE1	1.84	0.42
1:G:100:GLY:HA3	1:G:126:PHE:CD1	2.54	0.42
1:H:154:LEU:HB3	1:H:177:TYR:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122[A]:LEU:HB2	1:B:122:LEU:HD12	2.02	0.41
1:E:10:ARG:HH12	1:E:12:LYS:HE3	1.85	0.41
1:D:163:LYS:NZ	2:D:306:HOH:O	2.40	0.41
1:F:34:PRO:HA	1:F:69:PHE:HA	2.01	0.41
1:A:197:ARG:NH1	2:A:311:HOH:O	2.46	0.41
1:B:118:LYS:HB2	1:B:118:LYS:HE2	1.83	0.41
1:B:28:GLY:HA2	1:B:42:ARG:O	2.20	0.41
1:G:106:GLN:HA	1:G:118:LYS:O	2.21	0.41
1:C:205:LYS:HA	1:C:205:LYS:HD2	1.91	0.41
1:A:122[A]:LEU:HD23	1:A:122[A]:LEU:HA	1.87	0.41
1:B:138:GLU:OE2	2:B:302:HOH:O	2.21	0.40
1:B:19:VAL:HG23	1:B:24:PHE:HE1	1.86	0.40

All (1) 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:C:7:GLU:N	1:D:7:GLU:OE2[3_545]	2.10	0.10

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	221/228 (97%)	217 (98%)	4 (2%)	0	100	100
1	B	223/228 (98%)	219 (98%)	4 (2%)	0	100	100
1	C	220/228 (96%)	216 (98%)	4 (2%)	0	100	100
1	D	224/228 (98%)	220 (98%)	4 (2%)	0	100	100
1	E	222/228 (97%)	218 (98%)	4 (2%)	0	100	100
1	F	219/228 (96%)	215 (98%)	4 (2%)	0	100	100
1	G	218/228 (96%)	214 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	217/228 (95%)	213 (98%)	4 (2%)	0	100	100
All	All	1764/1824 (97%)	1732 (98%)	32 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	192/196 (98%)	190 (99%)	2 (1%)	76	75
1	B	194/196 (99%)	192 (99%)	2 (1%)	76	75
1	C	191/196 (97%)	189 (99%)	2 (1%)	76	75
1	D	195/196 (100%)	193 (99%)	2 (1%)	76	75
1	E	194/196 (99%)	192 (99%)	2 (1%)	76	75
1	F	191/196 (97%)	187 (98%)	4 (2%)	53	48
1	G	189/196 (96%)	187 (99%)	2 (1%)	73	72
1	H	189/196 (96%)	187 (99%)	2 (1%)	73	72
All	All	1535/1568 (98%)	1517 (99%)	18 (1%)	73	69

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ARG
1	A	67	ARG
1	B	42	ARG
1	B	131	PRO
1	C	42	ARG
1	C	113	ASN
1	D	12	LYS
1	D	42	ARG
1	E	42	ARG
1	E	120	LYS

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Mol	Chain	Res	Type
1	F	18[A]	THR
1	F	18[B]	THR
1	F	42	ARG
1	F	67	ARG
1	G	42	ARG
1	G	167	ARG
1	H	13	MET
1	H	138	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	105	HIS
1	F	44	HIS
1	F	72	HIS
1	G	44	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](#)

There are no non-standard protein/DNA/RNA residues in this entry.

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 ⓘ

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, 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	220/228 (96%)	-0.33	4 (1%) 68 70	14, 21, 41, 62	0
1	B	220/228 (96%)	-0.29	3 (1%) 75 76	11, 19, 42, 55	0
1	C	220/228 (96%)	-0.05	7 (3%) 47 50	15, 28, 53, 83	0
1	D	220/228 (96%)	-0.16	5 (2%) 60 63	12, 23, 45, 80	0
1	E	219/228 (96%)	0.15	10 (4%) 32 33	20, 35, 60, 77	0
1	F	219/228 (96%)	0.12	8 (3%) 41 43	25, 36, 59, 79	0
1	G	220/228 (96%)	0.61	29 (13%) 3 2	23, 46, 77, 99	0
1	H	219/228 (96%)	0.71	29 (13%) 3 2	28, 50, 75, 97	0
All	All	1757/1824 (96%)	0.10	95 (5%) 25 26	11, 31, 65, 99	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	4	LEU	6.9
1	H	208	TYR	5.6
1	F	202	ARG	5.5
1	C	183	ILE	5.3
1	C	182	ALA	5.1
1	F	225	LYS	5.0
1	H	113	ASN	5.0
1	H	184	ARG	4.8
1	H	112	GLY	4.7
1	G	184	ARG	4.6
1	H	183	ILE	4.6
1	G	5	LEU	4.4
1	G	225	LYS	4.4
1	H	205	LYS	4.3
1	C	184	ARG	4.3
1	E	166	ASP	4.1

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Mol	Chain	Res	Type	RSRZ
1	D	3	GLY	4.0
1	G	3	GLY	4.0
1	D	184	ARG	4.0
1	H	225	LYS	4.0
1	G	110	LEU	3.9
1	H	6	LYS	3.8
1	G	208	TYR	3.7
1	H	110	LEU	3.6
1	C	4	LEU	3.5
1	C	3	GLY	3.5
1	G	181	LYS	3.4
1	H	4	LEU	3.3
1	H	182	ALA	3.2
1	B	225	LYS	3.1
1	H	186	LEU	3.1
1	G	186	LEU	3.1
1	A	184	ARG	3.0
1	G	111	GLU	3.0
1	G	10	ARG	3.0
1	H	75	GLY	3.0
1	H	10	ARG	2.9
1	H	25	LYS	2.9
1	H	202	ARG	2.9
1	E	202	ARG	2.8
1	G	185	ALA	2.8
1	F	205	LYS	2.8
1	G	183	ILE	2.8
1	E	129	ASP	2.8
1	G	35	PHE	2.8
1	H	47	GLU	2.8
1	G	112	GLY	2.7
1	H	151	ASN	2.7
1	A	225	LYS	2.7
1	F	7	GLU	2.6
1	G	205	LYS	2.6
1	B	183	ILE	2.6
1	H	7	GLU	2.6
1	G	7	GLU	2.5
1	D	67[A]	ARG	2.5
1	H	204	LYS	2.5
1	F	204	LYS	2.5
1	G	182	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	G	202	ARG	2.5
1	E	184	ARG	2.4
1	E	7	GLU	2.4
1	G	114	CYS	2.4
1	H	116	ILE	2.4
1	A	3	GLY	2.4
1	D	183	ILE	2.4
1	E	167	ARG	2.4
1	F	208	TYR	2.4
1	G	86	GLU	2.3
1	H	185	ALA	2.3
1	F	184	ARG	2.3
1	H	44	HIS	2.3
1	A	183	ILE	2.3
1	E	44	HIS	2.2
1	F	167	ARG	2.2
1	E	205	LYS	2.2
1	G	6	LYS	2.2
1	H	46	THR	2.2
1	B	3	GLY	2.2
1	H	167	ARG	2.1
1	H	60	PRO	2.1
1	G	129	ASP	2.1
1	G	90	TRP	2.1
1	C	60	PRO	2.1
1	C	61	CYS	2.1
1	G	45	VAL	2.1
1	D	185	ALA	2.1
1	G	12	LYS	2.1
1	E	204	LYS	2.1
1	G	224	GLU	2.0
1	H	66	SER	2.0
1	G	60	PRO	2.0
1	E	183	ILE	2.0
1	H	111	GLU	2.0
1	H	21	GLY	2.0
1	G	67	ARG	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

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.