



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 02:03 PM UTC

PDB ID : 3VIC / pdb_00003vic
Title : Green-fluorescent variant of the non-fluorescent chromoprotein Rtms5
Authors : Battad, J.M.; Traore, D.A.K.; Byres, E.; Wilce, M.; Devenish, R.J.; Rossjohn, J.; Prescott, M.
Deposited on : 2011-09-28
Resolution : 2.20 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:

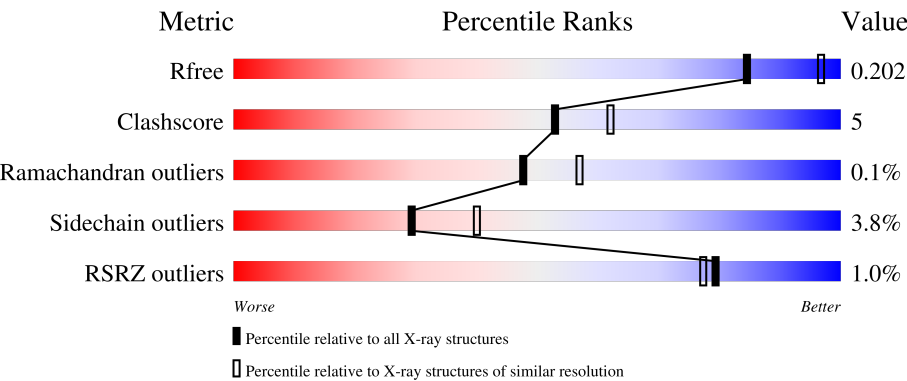
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	238	79%11%• 8%
1	B	238	% 79%11%• 8%
1	C	238	% 82%9%• 8%
1	D	238	% 77%13%• 8%
1	E	238	% 78%11%• 8%

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Mol	Chain	Length	Quality of chain
1	F	238	77% 13% 8%
1	G	238	77% 14% 8%
1	H	238	77% 13% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15565 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 GFP-like non-fluorescent chromoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	0	4	0
			1772	1134	295	330	13			
1	B	219	Total	C	N	O	S	0	3	0
			1766	1130	294	329	13			
1	C	219	Total	C	N	O	S	0	2	0
			1766	1130	294	329	13			
1	D	219	Total	C	N	O	S	0	3	0
			1766	1130	294	329	13			
1	E	218	Total	C	N	O	S	0	3	0
			1761	1127	293	328	13			
1	F	219	Total	C	N	O	S	0	2	0
			1762	1128	293	327	14			
1	G	219	Total	C	N	O	S	0	1	0
			1751	1122	291	325	13			
1	H	218	Total	C	N	O	S	0	2	0
			1754	1123	292	326	13			

There are 176 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MET	-	expression tag	UNP P83690
A	-13	ARG	-	expression tag	UNP P83690
A	-12	GLY	-	expression tag	UNP P83690
A	-11	SER	-	expression tag	UNP P83690
A	-10	HIS	-	expression tag	UNP P83690
A	-9	HIS	-	expression tag	UNP P83690
A	-8	HIS	-	expression tag	UNP P83690
A	-7	HIS	-	expression tag	UNP P83690
A	-6	HIS	-	expression tag	UNP P83690
A	-5	HIS	-	expression tag	UNP P83690
A	-4	THR	-	expression tag	UNP P83690
A	-3	ASP	-	expression tag	UNP P83690
A	-2	PRO	-	expression tag	UNP P83690

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP P83690
A	0	SER	-	expression tag	UNP P83690
A	1	ARG	-	expression tag	UNP P83690
A	2	SER	-	expression tag	UNP P83690
A	3	LYS	-	expression tag	UNP P83690
A	4	THR	-	expression tag	UNP P83690
A	66	QFG	GLN	chromophore	UNP P83690
A	66	QFG	TYR	chromophore	UNP P83690
A	66	QFG	GLY	chromophore	UNP P83690
B	-14	MET	-	expression tag	UNP P83690
B	-13	ARG	-	expression tag	UNP P83690
B	-12	GLY	-	expression tag	UNP P83690
B	-11	SER	-	expression tag	UNP P83690
B	-10	HIS	-	expression tag	UNP P83690
B	-9	HIS	-	expression tag	UNP P83690
B	-8	HIS	-	expression tag	UNP P83690
B	-7	HIS	-	expression tag	UNP P83690
B	-6	HIS	-	expression tag	UNP P83690
B	-5	HIS	-	expression tag	UNP P83690
B	-4	THR	-	expression tag	UNP P83690
B	-3	ASP	-	expression tag	UNP P83690
B	-2	PRO	-	expression tag	UNP P83690
B	-1	SER	-	expression tag	UNP P83690
B	0	SER	-	expression tag	UNP P83690
B	1	ARG	-	expression tag	UNP P83690
B	2	SER	-	expression tag	UNP P83690
B	3	LYS	-	expression tag	UNP P83690
B	4	THR	-	expression tag	UNP P83690
B	66	QFG	GLN	chromophore	UNP P83690
B	66	QFG	TYR	chromophore	UNP P83690
B	66	QFG	GLY	chromophore	UNP P83690
C	-14	MET	-	expression tag	UNP P83690
C	-13	ARG	-	expression tag	UNP P83690
C	-12	GLY	-	expression tag	UNP P83690
C	-11	SER	-	expression tag	UNP P83690
C	-10	HIS	-	expression tag	UNP P83690
C	-9	HIS	-	expression tag	UNP P83690
C	-8	HIS	-	expression tag	UNP P83690
C	-7	HIS	-	expression tag	UNP P83690
C	-6	HIS	-	expression tag	UNP P83690
C	-5	HIS	-	expression tag	UNP P83690
C	-4	THR	-	expression tag	UNP P83690

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	ASP	-	expression tag	UNP P83690
C	-2	PRO	-	expression tag	UNP P83690
C	-1	SER	-	expression tag	UNP P83690
C	0	SER	-	expression tag	UNP P83690
C	1	ARG	-	expression tag	UNP P83690
C	2	SER	-	expression tag	UNP P83690
C	3	LYS	-	expression tag	UNP P83690
C	4	THR	-	expression tag	UNP P83690
C	66	QFG	GLN	chromophore	UNP P83690
C	66	QFG	TYR	chromophore	UNP P83690
C	66	QFG	GLY	chromophore	UNP P83690
D	-14	MET	-	expression tag	UNP P83690
D	-13	ARG	-	expression tag	UNP P83690
D	-12	GLY	-	expression tag	UNP P83690
D	-11	SER	-	expression tag	UNP P83690
D	-10	HIS	-	expression tag	UNP P83690
D	-9	HIS	-	expression tag	UNP P83690
D	-8	HIS	-	expression tag	UNP P83690
D	-7	HIS	-	expression tag	UNP P83690
D	-6	HIS	-	expression tag	UNP P83690
D	-5	HIS	-	expression tag	UNP P83690
D	-4	THR	-	expression tag	UNP P83690
D	-3	ASP	-	expression tag	UNP P83690
D	-2	PRO	-	expression tag	UNP P83690
D	-1	SER	-	expression tag	UNP P83690
D	0	SER	-	expression tag	UNP P83690
D	1	ARG	-	expression tag	UNP P83690
D	2	SER	-	expression tag	UNP P83690
D	3	LYS	-	expression tag	UNP P83690
D	4	THR	-	expression tag	UNP P83690
D	66	QFG	GLN	chromophore	UNP P83690
D	66	QFG	TYR	chromophore	UNP P83690
D	66	QFG	GLY	chromophore	UNP P83690
E	-14	MET	-	expression tag	UNP P83690
E	-13	ARG	-	expression tag	UNP P83690
E	-12	GLY	-	expression tag	UNP P83690
E	-11	SER	-	expression tag	UNP P83690
E	-10	HIS	-	expression tag	UNP P83690
E	-9	HIS	-	expression tag	UNP P83690
E	-8	HIS	-	expression tag	UNP P83690
E	-7	HIS	-	expression tag	UNP P83690
E	-6	HIS	-	expression tag	UNP P83690

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-5	HIS	-	expression tag	UNP P83690
E	-4	THR	-	expression tag	UNP P83690
E	-3	ASP	-	expression tag	UNP P83690
E	-2	PRO	-	expression tag	UNP P83690
E	-1	SER	-	expression tag	UNP P83690
E	0	SER	-	expression tag	UNP P83690
E	1	ARG	-	expression tag	UNP P83690
E	2	SER	-	expression tag	UNP P83690
E	3	LYS	-	expression tag	UNP P83690
E	4	THR	-	expression tag	UNP P83690
E	66	QFG	GLN	chromophore	UNP P83690
E	66	QFG	TYR	chromophore	UNP P83690
E	66	QFG	GLY	chromophore	UNP P83690
F	-14	MET	-	expression tag	UNP P83690
F	-13	ARG	-	expression tag	UNP P83690
F	-12	GLY	-	expression tag	UNP P83690
F	-11	SER	-	expression tag	UNP P83690
F	-10	HIS	-	expression tag	UNP P83690
F	-9	HIS	-	expression tag	UNP P83690
F	-8	HIS	-	expression tag	UNP P83690
F	-7	HIS	-	expression tag	UNP P83690
F	-6	HIS	-	expression tag	UNP P83690
F	-5	HIS	-	expression tag	UNP P83690
F	-4	THR	-	expression tag	UNP P83690
F	-3	ASP	-	expression tag	UNP P83690
F	-2	PRO	-	expression tag	UNP P83690
F	-1	SER	-	expression tag	UNP P83690
F	0	SER	-	expression tag	UNP P83690
F	1	ARG	-	expression tag	UNP P83690
F	2	SER	-	expression tag	UNP P83690
F	3	LYS	-	expression tag	UNP P83690
F	4	THR	-	expression tag	UNP P83690
F	66	QFG	GLN	chromophore	UNP P83690
F	66	QFG	TYR	chromophore	UNP P83690
F	66	QFG	GLY	chromophore	UNP P83690
G	-14	MET	-	expression tag	UNP P83690
G	-13	ARG	-	expression tag	UNP P83690
G	-12	GLY	-	expression tag	UNP P83690
G	-11	SER	-	expression tag	UNP P83690
G	-10	HIS	-	expression tag	UNP P83690
G	-9	HIS	-	expression tag	UNP P83690
G	-8	HIS	-	expression tag	UNP P83690

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-7	HIS	-	expression tag	UNP P83690
G	-6	HIS	-	expression tag	UNP P83690
G	-5	HIS	-	expression tag	UNP P83690
G	-4	THR	-	expression tag	UNP P83690
G	-3	ASP	-	expression tag	UNP P83690
G	-2	PRO	-	expression tag	UNP P83690
G	-1	SER	-	expression tag	UNP P83690
G	0	SER	-	expression tag	UNP P83690
G	1	ARG	-	expression tag	UNP P83690
G	2	SER	-	expression tag	UNP P83690
G	3	LYS	-	expression tag	UNP P83690
G	4	THR	-	expression tag	UNP P83690
G	66	QFG	GLN	chromophore	UNP P83690
G	66	QFG	TYR	chromophore	UNP P83690
G	66	QFG	GLY	chromophore	UNP P83690
H	-14	MET	-	expression tag	UNP P83690
H	-13	ARG	-	expression tag	UNP P83690
H	-12	GLY	-	expression tag	UNP P83690
H	-11	SER	-	expression tag	UNP P83690
H	-10	HIS	-	expression tag	UNP P83690
H	-9	HIS	-	expression tag	UNP P83690
H	-8	HIS	-	expression tag	UNP P83690
H	-7	HIS	-	expression tag	UNP P83690
H	-6	HIS	-	expression tag	UNP P83690
H	-5	HIS	-	expression tag	UNP P83690
H	-4	THR	-	expression tag	UNP P83690
H	-3	ASP	-	expression tag	UNP P83690
H	-2	PRO	-	expression tag	UNP P83690
H	-1	SER	-	expression tag	UNP P83690
H	0	SER	-	expression tag	UNP P83690
H	1	ARG	-	expression tag	UNP P83690
H	2	SER	-	expression tag	UNP P83690
H	3	LYS	-	expression tag	UNP P83690
H	4	THR	-	expression tag	UNP P83690
H	66	QFG	GLN	chromophore	UNP P83690
H	66	QFG	TYR	chromophore	UNP P83690
H	66	QFG	GLY	chromophore	UNP P83690

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total I 1 1	0	0
2	B	1	Total I 1 1	0	0
2	C	2	Total I 2 2	0	0
2	D	2	Total I 2 2	0	0
2	E	1	Total I 1 1	0	0
2	H	1	Total I 1 1	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Cl 2 2	0	0
3	B	2	Total Cl 2 2	0	0
3	C	2	Total Cl 2 2	0	0
3	D	2	Total Cl 2 2	0	0
3	E	2	Total Cl 2 2	0	0
3	F	2	Total Cl 2 2	0	0
3	G	2	Total Cl 2 2	0	0
3	H	2	Total Cl 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	184	Total O 184 184	0	0
4	B	188	Total O 188 188	0	0

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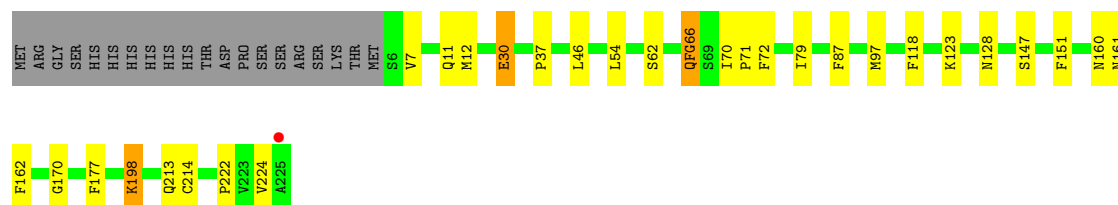
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	169	Total 169	O 169	0	0
4	D	185	Total 185	O 185	0	0
4	E	193	Total 193	O 193	0	0
4	F	189	Total 189	O 189	0	0
4	G	179	Total 179	O 179	0	0
4	H	156	Total 156	O 156	0	0

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.

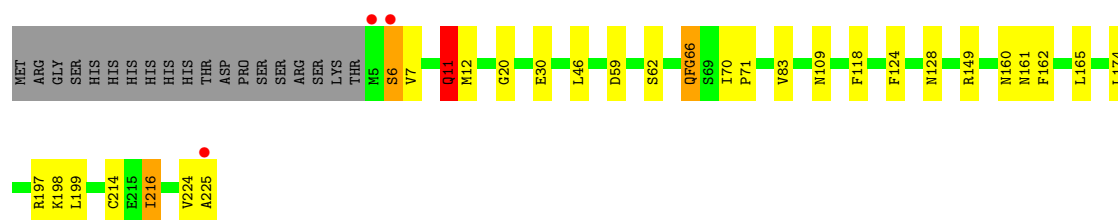
- Molecule 1: GFP-like non-fluorescent chromoprotein

Chain A: 



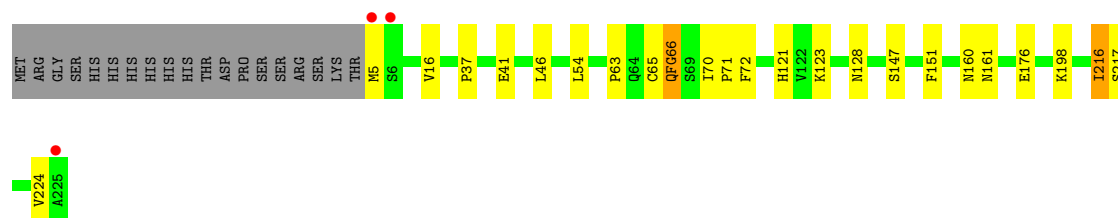
- Molecule 1: GFP-like non-fluorescent chromoprotein

Chain B: 



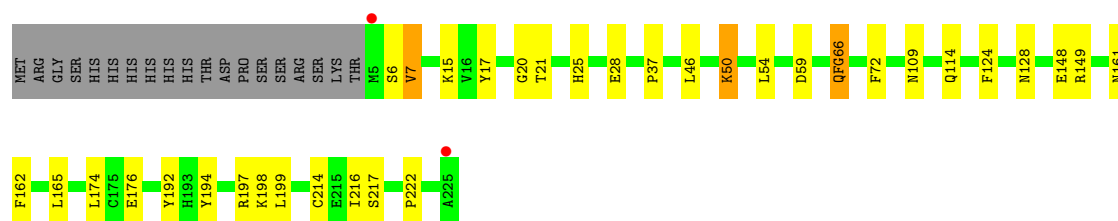
- Molecule 1: GFP-like non-fluorescent chromoprotein

Chain C: 

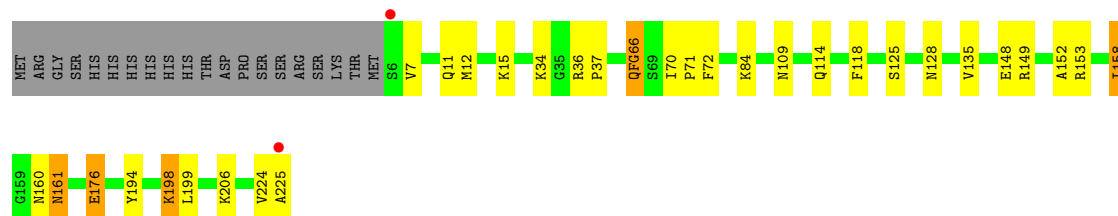
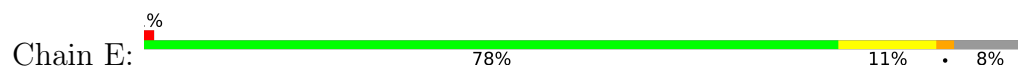


- Molecule 1: GFP-like non-fluorescent chromoprotein

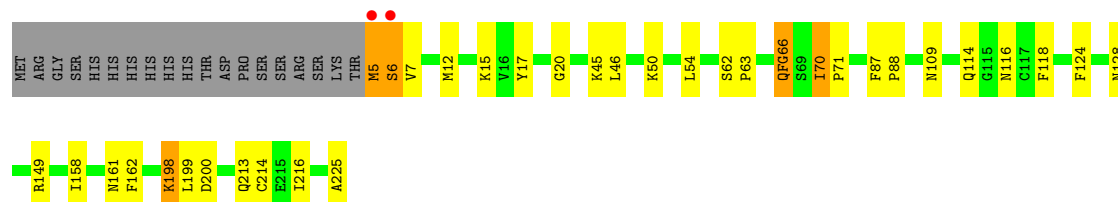
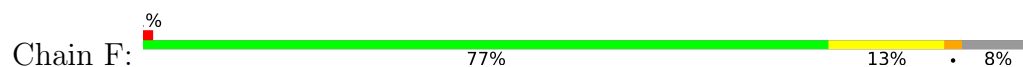
Chain D: 



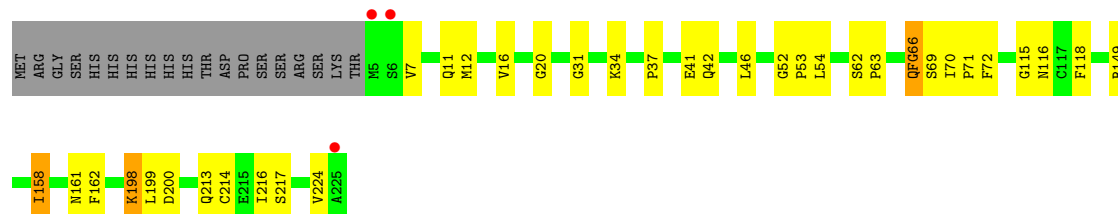
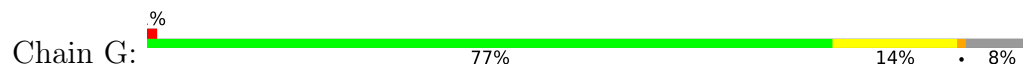
- Molecule 1: GFP-like non-fluorescent chromoprotein



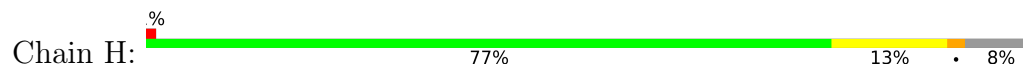
- Molecule 1: GFP-like non-fluorescent chromoprotein

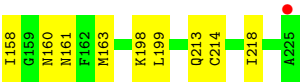


- Molecule 1: GFP-like non-fluorescent chromoprotein



- Molecule 1: GFP-like non-fluorescent chromoprotein





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	150.27Å 186.06Å 185.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.80 – 2.20 54.80 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (54.80-2.20) 100.0 (54.80-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.48 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.160 , 0.204 0.158 , 0.202	Depositor DCC
R_{free} test set	6587 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15565	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.99% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, QFG, IOD

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	1.24	2/1798 (0.1%)	1.10	7/2432 (0.3%)
1	B	1.26	1/1795 (0.1%)	1.10	6/2428 (0.2%)
1	C	1.24	2/1792 (0.1%)	1.04	1/2423 (0.0%)
1	D	1.29	2/1795 (0.1%)	1.11	1/2428 (0.0%)
1	E	1.31	4/1790 (0.2%)	1.11	6/2421 (0.2%)
1	F	1.27	2/1791 (0.1%)	1.07	3/2421 (0.1%)
1	G	1.21	1/1780 (0.1%)	1.07	2/2407 (0.1%)
1	H	1.25	1/1783 (0.1%)	1.10	6/2411 (0.2%)
All	All	1.26	15/14324 (0.1%)	1.09	32/19371 (0.2%)

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

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	158	ILE	CA-CB	7.25	1.62	1.54
1	A	222	PRO	CA-C	5.92	1.59	1.52
1	H	218	ILE	CA-CB	5.81	1.61	1.54
1	D	222	PRO	N-CA	5.59	1.53	1.47
1	E	135	VAL	CA-CB	5.48	1.60	1.54
1	E	158	ILE	CA-C	-5.44	1.46	1.52
1	F	6	SER	C-O	5.33	1.31	1.24
1	E	176	GLU	CA-C	5.25	1.59	1.52
1	E	152	ALA	CA-CB	5.17	1.61	1.53
1	D	7	VAL	C-O	5.14	1.31	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	123	LYS	N-CA	5.11	1.52	1.46
1	C	216	ILE	CA-CB	5.11	1.60	1.54
1	A	30	GLU	CA-C	-5.06	1.46	1.52
1	F	50	LYS	C-O	-5.01	1.18	1.23
1	B	11	GLN	CG-CD	5.00	1.64	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	62	SER	CA-C-N	-7.04	112.63	120.45
1	B	62	SER	C-N-CA	-7.04	112.63	120.45
1	B	161	ASN	CB-CA-C	-6.34	96.79	109.35
1	A	161	ASN	CB-CA-C	-6.14	96.91	109.38
1	D	161	ASN	CB-CA-C	-6.11	96.97	109.38
1	H	7	VAL	N-CA-C	-6.09	104.43	113.39
1	A	7	VAL	N-CA-C	6.07	116.81	110.62
1	A	62	SER	CA-C-N	-6.04	113.58	119.87
1	A	62	SER	C-N-CA	-6.04	113.58	119.87
1	H	158	ILE	CG1-CB-CG2	-5.96	92.81	110.70
1	A	79	ILE	CA-C-N	-5.77	114.01	119.90
1	A	79	ILE	C-N-CA	-5.77	114.01	119.90
1	H	161	ASN	CB-CA-C	-5.74	97.73	109.38
1	F	161	ASN	CB-CA-C	-5.58	98.31	109.35
1	G	20	GLY	N-CA-C	5.56	117.68	110.45
1	E	161	ASN	N-CA-CB	5.46	120.90	111.13
1	H	20	GLY	N-CA-C	5.37	118.09	110.74
1	B	6	SER	N-CA-C	-5.34	105.54	111.36
1	E	125	SER	CA-C-N	5.32	125.27	121.65
1	E	125	SER	C-N-CA	5.32	125.27	121.65
1	B	83	VAL	N-CA-C	5.31	115.52	110.42
1	E	36	ARG	CA-C-N	5.29	124.92	119.05
1	E	36	ARG	C-N-CA	5.29	124.92	119.05
1	E	161	ASN	CB-CA-C	-5.28	98.66	109.38
1	G	161	ASN	CB-CA-C	-5.28	98.66	109.38
1	C	161	ASN	CB-CA-C	-5.28	98.90	109.35
1	F	70	ILE	CA-C-N	-5.21	114.45	119.87
1	F	70	ILE	C-N-CA	-5.21	114.45	119.87
1	H	161	ASN	N-CA-CB	5.21	120.46	111.13
1	H	16	VAL	CB-CA-C	-5.15	103.16	110.83
1	B	7	VAL	CB-CA-C	-5.02	105.36	112.14
1	A	87	PHE	CB-CA-C	5.01	116.63	109.11

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	6	SER	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	1772	0	1702	19	0
1	B	1766	0	1695	20	0
1	C	1766	0	1697	15	0
1	D	1766	0	1695	20	0
1	E	1761	0	1693	20	0
1	F	1762	0	1696	24	0
1	G	1751	0	1684	20	0
1	H	1754	0	1687	17	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	1	0	0	0	0
2	H	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
3	G	2	0	0	0	0
3	H	2	0	0	0	0
4	A	184	0	0	7	0
4	B	188	0	0	6	0
4	C	169	0	0	5	0
4	D	185	0	0	6	0
4	E	193	0	0	7	0
4	F	189	0	0	6	0
4	G	179	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	156	0	0	5	0
All	All	15565	0	13549	137	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 (137) 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:66:QFG:HB2	4:A:1086:HOH:O	1.25	1.35
1:F:66:QFG:HB2	4:F:1218:HOH:O	1.21	1.35
1:D:66:QFG:HB2	4:D:260:HOH:O	1.22	1.34
1:C:66:QFG:HB2	4:C:1217:HOH:O	1.24	1.31
1:G:66:QFG:HB2	4:G:1219:HOH:O	1.25	1.30
1:B:66:QFG:HB2	4:B:1214:HOH:O	1.18	1.30
1:E:66:QFG:HB2	4:E:1146:HOH:O	1.30	1.29
1:D:128[B]:ASN:OD1	4:D:347:HOH:O	1.62	1.15
1:H:66:QFG:HB2	4:H:1220:HOH:O	0.98	1.15
1:F:128[A]:ASN:OD1	4:F:1297:HOH:O	1.69	1.09
1:A:128[B]:ASN:OD1	4:A:466:HOH:O	1.69	1.08
1:A:11:GLN:NE2	4:A:1081:HOH:O	1.97	0.96
1:B:214[B]:CYS:SG	4:B:1252:HOH:O	2.27	0.91
1:G:214[B]:CYS:SG	4:G:1349:HOH:O	2.36	0.83
1:C:176:GLU:OE2	4:C:1439:HOH:O	2.00	0.79
1:G:12:MET:HE2	1:G:118:PHE:CE2	2.19	0.78
1:F:128[A]:ASN:ND2	4:F:1296:HOH:O	2.06	0.76
1:B:11:GLN:OE1	4:B:1100:HOH:O	2.07	0.72
1:G:214[A]:CYS:SG	4:G:1360:HOH:O	2.48	0.72
1:H:108:SER:OG	4:H:1310:HOH:O	2.08	0.70
1:C:54:LEU:HD12	4:C:1255:HOH:O	1.94	0.68
1:F:12:MET:HE2	1:F:118:PHE:CE2	2.30	0.66
1:A:170:GLY:O	4:A:1242:HOH:O	2.14	0.66
1:F:5:MET:HB3	1:F:7:VAL:H	1.61	0.64
1:H:12:MET:HE2	1:H:118:PHE:CE2	2.37	0.60
1:E:37:PRO:HA	1:E:72:PHE:HA	1.85	0.59
1:E:198:LYS:HD3	4:E:393:HOH:O	2.03	0.58
1:A:12:MET:HE2	1:A:118:PHE:CE2	2.38	0.57
1:A:54:LEU:HD12	4:A:973:HOH:O	2.04	0.57
1:G:54:LEU:HD12	4:G:1403:HOH:O	2.05	0.56
1:E:128[B]:ASN:ND2	4:E:380:HOH:O	2.29	0.56
1:E:198:LYS:NZ	1:F:225:ALA:C	2.64	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:LYS:HZ3	1:B:225:ALA:C	2.14	0.55
1:C:151:PHE:CZ	1:D:174:LEU:HD11	2.42	0.55
1:D:214[B]:CYS:SG	4:D:899:HOH:O	2.58	0.55
1:A:12:MET:HE2	1:A:118:PHE:CZ	2.41	0.55
1:E:198:LYS:HZ1	1:F:225:ALA:C	2.15	0.55
1:A:30:GLU:OE2	4:A:1432:HOH:O	2.18	0.54
1:B:12:MET:HE3	1:B:118:PHE:CE2	2.42	0.54
1:F:214[A]:CYS:SG	4:F:1390:HOH:O	2.32	0.54
1:F:54:LEU:HD12	4:F:886:HOH:O	2.07	0.53
1:H:15:LYS:HG2	1:H:17:TYR:CE1	2.44	0.53
1:C:128[B]:ASN:ND2	4:C:664:HOH:O	2.30	0.52
1:E:160:ASN:HB3	1:F:162:PHE:CZ	2.44	0.52
1:H:12:MET:HE2	1:H:118:PHE:CZ	2.44	0.52
1:G:224:VAL:HG12	1:H:198:LYS:HG2	1.92	0.52
1:A:198:LYS:NZ	1:B:225:ALA:C	2.68	0.52
1:A:37:PRO:HA	1:A:72:PHE:HA	1.93	0.51
1:D:54:LEU:HD12	4:D:1257:HOH:O	2.10	0.51
1:F:70:ILE:N	1:F:71:PRO:CD	2.73	0.51
1:G:12:MET:HE2	1:G:118:PHE:CZ	2.45	0.51
1:E:206:LYS:NZ	4:E:1085:HOH:O	2.43	0.50
1:G:7:VAL:O	4:G:1340:HOH:O	2.18	0.50
1:G:162:PHE:CZ	1:H:160:ASN:HB3	2.46	0.50
1:E:224:VAL:HG12	1:F:198:LYS:HB2	1.94	0.50
1:E:12:MET:HE2	1:E:118:PHE:CZ	2.47	0.49
1:G:198:LYS:HE3	1:G:200:ASP:OD2	2.11	0.49
1:F:116:ASN:HB3	4:F:1417:HOH:O	2.11	0.49
1:G:116:ASN:OD1	4:G:1341:HOH:O	2.20	0.49
1:A:162:PHE:CZ	1:B:160:ASN:HB3	2.48	0.48
1:E:12:MET:HE3	1:E:118:PHE:CE2	2.48	0.48
1:B:70:ILE:N	1:B:71:PRO:CD	2.76	0.48
1:E:225:ALA:OXT	1:F:198:LYS:HD3	2.12	0.48
1:G:70:ILE:N	1:G:71:PRO:CD	2.76	0.48
1:B:20:GLY:HA3	1:B:124:PHE:O	2.13	0.48
1:E:12:MET:CE	1:E:118:PHE:CE2	2.96	0.48
1:D:162:PHE:HE2	1:D:176:GLU:OE2	1.97	0.48
1:E:11:GLN:NE2	4:E:1103:HOH:O	2.23	0.47
1:H:42:GLN:HE22	1:H:69:SER:HB3	1.80	0.47
1:E:176:GLU:OE1	4:E:1267:HOH:O	2.20	0.47
1:A:224:VAL:HG12	1:B:198:LYS:HG2	1.95	0.46
1:C:65:CYS:O	1:C:66:QFG:HG1	2.15	0.46
1:E:70:ILE:N	1:E:71:PRO:CD	2.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:63:PRO:O	1:G:66:QFG:C2	2.64	0.46
1:H:214[B]:CYS:SG	4:H:1338:HOH:O	2.61	0.46
1:C:224:VAL:HG12	1:D:198:LYS:HG2	1.97	0.46
1:H:54:LEU:HD12	4:H:496:HOH:O	2.15	0.46
1:F:12:MET:HE2	1:F:118:PHE:CZ	2.51	0.46
1:H:70:ILE:N	1:H:71:PRO:CD	2.79	0.46
1:C:41:GLU:HG2	1:C:216:ILE:HD12	1.98	0.45
1:E:153:ARG:HB3	1:E:158:ILE:CD1	2.46	0.45
1:E:148:GLU:O	1:E:194:TYR:HA	2.16	0.45
1:C:121:HIS:HD2	4:C:1018:HOH:O	2.00	0.45
1:H:37:PRO:HA	1:H:72:PHE:HA	1.98	0.45
1:D:176:GLU:HG2	4:D:394:HOH:O	2.17	0.45
1:F:198:LYS:HE3	1:F:198:LYS:HB3	1.62	0.45
1:A:151:PHE:CZ	1:B:174:LEU:HD11	2.52	0.45
1:A:70:ILE:N	1:A:71:PRO:CD	2.80	0.45
1:H:128[B]:ASN:ND2	4:H:342:HOH:O	2.33	0.44
1:G:37:PRO:HA	1:G:72:PHE:HA	1.99	0.44
1:B:160:ASN:ND2	4:B:1264:HOH:O	2.14	0.44
1:D:197:ARG:HA	1:D:216:ILE:O	2.16	0.44
1:E:84:LYS:NZ	4:E:239:HOH:O	2.51	0.44
1:A:123:LYS:NZ	4:A:611:HOH:O	2.50	0.44
1:C:70:ILE:N	1:C:71:PRO:CD	2.81	0.44
1:F:198:LYS:HD2	1:F:200:ASP:OD2	2.18	0.44
1:A:160:ASN:HB3	1:B:162:PHE:CZ	2.52	0.43
1:B:30:GLU:OE2	4:B:1247:HOH:O	2.21	0.43
1:B:128[B]:ASN:ND2	4:B:747:HOH:O	2.30	0.43
1:H:146:HIS:CG	1:H:163:MET:SD	3.11	0.43
1:C:37:PRO:HA	1:C:72:PHE:HA	2.00	0.43
1:C:198:LYS:HE3	1:C:216:ILE:HG12	2.01	0.43
1:H:63:PRO:O	1:H:66:QFG:C2	2.67	0.43
1:D:149:ARG:HD3	1:D:192:TYR:CZ	2.54	0.43
1:D:21:THR:HA	1:D:25:HIS:O	2.19	0.42
1:G:62:SER:HB2	1:G:213:GLN:OE1	2.18	0.42
1:B:59:ASP:HB3	1:B:165:LEU:HD21	2.00	0.42
1:D:20:GLY:HA3	1:D:124:PHE:O	2.19	0.42
1:D:37:PRO:HA	1:D:72:PHE:HA	2.00	0.42
1:D:28:GLU:HB2	1:D:50:LYS:HB2	2.01	0.42
1:A:198:LYS:HB3	1:B:224:VAL:HG12	2.01	0.42
1:C:160:ASN:HB3	1:D:162:PHE:CZ	2.54	0.42
1:F:62:SER:HB2	1:F:213:GLN:OE1	2.19	0.42
1:G:31:GLY:HA3	1:G:46:LEU:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128[B]:ASN:HD22	1:B:128[B]:ASN:HA	1.73	0.42
1:F:66:QFG:O2	1:F:66:QFG:CD2	2.68	0.42
1:D:15:LYS:HG2	1:D:17:TYR:CE1	2.55	0.41
1:F:20:GLY:HA3	1:F:124:PHE:O	2.20	0.41
1:H:12:MET:HG3	1:H:37:PRO:HG3	2.01	0.41
1:G:149:ARG:HD2	1:G:149:ARG:C	2.45	0.41
1:A:97:MET:HE3	1:A:177:PHE:CZ	2.56	0.41
1:D:59:ASP:HB3	1:D:165:LEU:HD21	2.03	0.41
1:F:149:ARG:HD2	1:F:149:ARG:C	2.45	0.41
1:G:42:GLN:HE22	1:G:69:SER:HB3	1.85	0.41
1:C:151:PHE:HZ	1:D:174:LEU:HD11	1.83	0.41
1:E:149:ARG:HD2	1:E:149:ARG:C	2.44	0.41
1:C:63:PRO:O	1:C:66:QFG:C2	2.68	0.41
1:D:176:GLU:CG	4:D:394:HOH:O	2.68	0.41
1:F:87:PHE:HB3	1:F:88:PRO:HA	2.02	0.40
1:D:148:GLU:O	1:D:194:TYR:HA	2.22	0.40
1:F:15:LYS:HG2	1:F:17:TYR:CE1	2.56	0.40
1:H:21:THR:HA	1:H:25:HIS:O	2.22	0.40
1:G:52:GLY:HA2	1:G:53:PRO:C	2.46	0.40
1:B:149:ARG:HD2	1:B:149:ARG:C	2.47	0.40
1:B:197:ARG:HA	1:B:216:ILE:O	2.22	0.40
1:F:62:SER:N	1:F:63:PRO:CD	2.85	0.40
1:G:41:GLU:HG2	1:G:216:ILE:HD13	2.04	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	217/238 (91%)	213 (98%)	4 (2%)	0	100	100
1	B	217/238 (91%)	214 (99%)	3 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	216/238 (91%)	213 (99%)	3 (1%)	0	100	100
1	D	217/238 (91%)	215 (99%)	2 (1%)	0	100	100
1	E	216/238 (91%)	211 (98%)	5 (2%)	0	100	100
1	F	216/238 (91%)	212 (98%)	4 (2%)	0	100	100
1	G	215/238 (90%)	210 (98%)	4 (2%)	1 (0%)	24	27
1	H	215/238 (90%)	213 (99%)	2 (1%)	0	100	100
All	All	1729/1904 (91%)	1701 (98%)	27 (2%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	115	GLY

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	192/207 (93%)	186 (97%)	6 (3%)	35	48
1	B	191/207 (92%)	185 (97%)	6 (3%)	35	48
1	C	191/207 (92%)	186 (97%)	5 (3%)	40	55
1	D	191/207 (92%)	184 (96%)	7 (4%)	30	41
1	E	191/207 (92%)	183 (96%)	8 (4%)	26	36
1	F	191/207 (92%)	181 (95%)	10 (5%)	21	26
1	G	189/207 (91%)	182 (96%)	7 (4%)	30	41
1	H	190/207 (92%)	181 (95%)	9 (5%)	23	31
All	All	1526/1656 (92%)	1468 (96%)	58 (4%)	29	40

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU

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Mol	Chain	Res	Type
1	A	147	SER
1	A	198	LYS
1	A	213	GLN
1	A	214[A]	CYS
1	A	214[B]	CYS
1	B	6	SER
1	B	11	GLN
1	B	46	LEU
1	B	109	ASN
1	B	199	LEU
1	B	216	ILE
1	C	5	MET
1	C	16	VAL
1	C	46	LEU
1	C	147	SER
1	C	217	SER
1	D	7	VAL
1	D	46	LEU
1	D	50	LYS
1	D	109	ASN
1	D	114	GLN
1	D	199	LEU
1	D	217	SER
1	E	7	VAL
1	E	15	LYS
1	E	34	LYS
1	E	109	ASN
1	E	114	GLN
1	E	161	ASN
1	E	198	LYS
1	E	199	LEU
1	F	5	MET
1	F	6	SER
1	F	45	LYS
1	F	46	LEU
1	F	109	ASN
1	F	114	GLN
1	F	158	ILE
1	F	198	LYS
1	F	199	LEU
1	F	216	ILE
1	G	11	GLN

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Mol	Chain	Res	Type
1	G	16	VAL
1	G	34	LYS
1	G	158	ILE
1	G	198	LYS
1	G	199	LEU
1	G	217	SER
1	H	6	SER
1	H	7	VAL
1	H	16	VAL
1	H	46	LEU
1	H	108	SER
1	H	127	LEU
1	H	147	SER
1	H	199	LEU
1	H	213	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	HIS
1	B	11	GLN
1	B	109	ASN
1	B	121	HIS
1	B	161	ASN
1	C	11	GLN
1	C	132	ASN
1	D	121	HIS
1	D	132	ASN
1	F	109	ASN
1	F	121	HIS
1	F	161	ASN
1	G	132	ASN
1	H	42	GLN
1	H	116	ASN
1	H	137	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 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	QFG	C	66	1	24,24,25	2.76	6 (25%)	26,32,34	3.80	10 (38%)
1	QFG	E	66	1	24,24,25	3.31	8 (33%)	26,32,34	3.89	7 (26%)
1	QFG	B	66	1	24,24,25	2.58	7 (29%)	26,32,34	3.30	6 (23%)
1	QFG	A	66	1	24,24,25	2.98	7 (29%)	26,32,34	3.38	6 (23%)
1	QFG	D	66	1	24,24,25	2.98	6 (25%)	26,32,34	3.35	9 (34%)
1	QFG	F	66	1	24,24,25	2.98	8 (33%)	26,32,34	3.58	6 (23%)
1	QFG	G	66	1	24,24,25	2.78	10 (41%)	26,32,34	4.19	7 (26%)
1	QFG	H	66	1	24,24,25	3.25	8 (33%)	26,32,34	4.17	6 (23%)

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	QFG	C	66	1	-	4/10/32/33	0/2/2/2
1	QFG	E	66	1	-	4/10/32/33	0/2/2/2
1	QFG	B	66	1	-	5/10/32/33	0/2/2/2
1	QFG	A	66	1	-	4/10/32/33	0/2/2/2
1	QFG	D	66	1	-	5/10/32/33	0/2/2/2
1	QFG	F	66	1	-	4/10/32/33	0/2/2/2
1	QFG	G	66	1	-	3/10/32/33	0/2/2/2
1	QFG	H	66	1	-	6/10/32/33	0/2/2/2

All (60) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	66	QFG	CB2-CA2	13.88	1.48	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	66	QFG	CB2-CA2	12.87	1.47	1.35
1	F	66	QFG	CB2-CA2	11.92	1.46	1.35
1	D	66	QFG	CB2-CA2	11.22	1.46	1.35
1	C	66	QFG	CB2-CA2	10.85	1.45	1.35
1	B	66	QFG	CB2-CA2	10.38	1.45	1.35
1	A	66	QFG	CB2-CA2	10.21	1.45	1.35
1	G	66	QFG	CB2-CA2	9.79	1.44	1.35
1	A	66	QFG	C1-CA1	5.32	1.55	1.48
1	A	66	QFG	C1-N2	4.76	1.43	1.33
1	D	66	QFG	O2-C2	4.45	1.32	1.23
1	A	66	QFG	O2-C2	4.26	1.31	1.23
1	E	66	QFG	C1-N2	4.20	1.42	1.33
1	C	66	QFG	C1-N2	4.05	1.41	1.33
1	D	66	QFG	C1-CA1	3.88	1.53	1.48
1	H	66	QFG	C1-N2	3.76	1.41	1.33
1	D	66	QFG	C1-N2	3.67	1.41	1.33
1	G	66	QFG	C1-N2	3.66	1.41	1.33
1	H	66	QFG	C1-CA1	3.60	1.53	1.48
1	H	66	QFG	O2-C2	3.48	1.30	1.23
1	G	66	QFG	C1-N3	3.28	1.43	1.38
1	H	66	QFG	CE2-CD2	3.25	1.44	1.38
1	D	66	QFG	C1-N3	3.15	1.43	1.38
1	F	66	QFG	C1-N2	3.12	1.39	1.33
1	G	66	QFG	CG2-CB2	3.10	1.52	1.46
1	C	66	QFG	C1-CA1	3.09	1.52	1.48
1	B	66	QFG	C2-N3	-3.08	1.32	1.40
1	F	66	QFG	CE2-CD2	3.04	1.44	1.38
1	F	66	QFG	C1-N3	2.90	1.42	1.38
1	F	66	QFG	CG2-CB2	2.81	1.52	1.46
1	A	66	QFG	C1-N3	2.78	1.42	1.38
1	G	66	QFG	C1-CA1	2.74	1.52	1.48
1	G	66	QFG	CE2-CD2	2.72	1.43	1.38
1	H	66	QFG	CA2-N2	-2.70	1.32	1.38
1	E	66	QFG	C1-N3	2.66	1.42	1.38
1	E	66	QFG	CG2-CB2	2.51	1.51	1.46
1	H	66	QFG	C1-N3	2.50	1.42	1.38
1	F	66	QFG	O2-C2	2.48	1.28	1.23
1	G	66	QFG	C2-N3	-2.45	1.34	1.40
1	G	66	QFG	O2-C2	2.45	1.28	1.23
1	A	66	QFG	CE2-CD2	2.44	1.43	1.38
1	D	66	QFG	CG2-CB2	2.40	1.51	1.46
1	C	66	QFG	C2-N3	-2.36	1.34	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	66	QFG	CE2-CZ	2.32	1.43	1.38
1	E	66	QFG	CE1-CD1	2.29	1.42	1.38
1	G	66	QFG	CE2-CZ	2.28	1.43	1.38
1	H	66	QFG	CB1-CA1	2.27	1.56	1.50
1	C	66	QFG	O2-C2	2.26	1.27	1.23
1	F	66	QFG	C2-N3	-2.26	1.34	1.40
1	C	66	QFG	C1-N3	2.25	1.41	1.38
1	A	66	QFG	C2-N3	-2.23	1.34	1.40
1	E	66	QFG	O2-C2	2.22	1.27	1.23
1	E	66	QFG	CE2-CD2	2.22	1.42	1.38
1	B	66	QFG	CE2-CD2	2.15	1.42	1.38
1	F	66	QFG	CB1-CA1	2.13	1.56	1.50
1	E	66	QFG	CD3-NE1	2.11	1.39	1.32
1	B	66	QFG	C1-N3	2.09	1.41	1.38
1	G	66	QFG	CD2-CG2	2.09	1.43	1.39
1	B	66	QFG	O2-C2	2.05	1.27	1.23
1	B	66	QFG	CA2-C2	-2.03	1.46	1.48

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	66	QFG	O2-C2-CA2	-16.85	120.27	131.02
1	G	66	QFG	O2-C2-CA2	-13.65	122.31	131.02
1	E	66	QFG	CG2-CB2-CA2	-10.30	117.62	129.87
1	F	66	QFG	CG2-CB2-CA2	-10.26	117.67	129.87
1	A	66	QFG	O2-C2-CA2	-10.24	124.49	131.02
1	F	66	QFG	O2-C2-CA2	-10.22	124.50	131.02
1	E	66	QFG	CA2-C2-N3	9.48	111.46	103.50
1	B	66	QFG	O2-C2-CA2	-9.29	125.09	131.02
1	C	66	QFG	CA2-C2-N3	8.86	110.94	103.50
1	D	66	QFG	CG2-CB2-CA2	-8.77	119.44	129.87
1	C	66	QFG	O2-C2-CA2	-8.71	125.46	131.02
1	C	66	QFG	CB2-CA2-C2	8.63	132.81	122.36
1	G	66	QFG	CA2-C2-N3	8.19	110.38	103.50
1	D	66	QFG	CB2-CA2-C2	8.12	132.20	122.36
1	E	66	QFG	CB2-CA2-C2	8.04	132.11	122.36
1	G	66	QFG	CB2-CA2-C2	7.81	131.82	122.36
1	G	66	QFG	CG2-CB2-CA2	-7.70	120.72	129.87
1	E	66	QFG	O2-C2-CA2	-7.67	126.12	131.02
1	A	66	QFG	CA2-C2-N3	7.59	109.87	103.50
1	F	66	QFG	CA2-C2-N3	7.58	109.87	103.50
1	C	66	QFG	CG2-CB2-CA2	-7.29	121.20	129.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	66	QFG	CB2-CA2-N2	-7.24	118.93	128.76
1	B	66	QFG	CB2-CA2-C2	7.22	131.10	122.36
1	H	66	QFG	CA2-C2-N3	6.93	109.32	103.50
1	B	66	QFG	CA2-C2-N3	6.83	109.24	103.50
1	C	66	QFG	CB2-CA2-N2	-6.56	119.85	128.76
1	G	66	QFG	CB2-CA2-N2	-6.53	119.90	128.76
1	B	66	QFG	CG2-CB2-CA2	-6.52	122.11	129.87
1	H	66	QFG	CG2-CB2-CA2	-6.44	122.21	129.87
1	A	66	QFG	CG2-CB2-CA2	-6.41	122.24	129.87
1	A	66	QFG	CB2-CA2-C2	5.87	129.48	122.36
1	B	66	QFG	CB2-CA2-N2	-5.75	120.95	128.76
1	D	66	QFG	CA2-C2-N3	5.68	108.27	103.50
1	H	66	QFG	CB2-CA2-C2	5.60	129.14	122.36
1	D	66	QFG	O2-C2-CA2	-5.45	127.55	131.02
1	F	66	QFG	CB2-CA2-C2	5.39	128.90	122.36
1	E	66	QFG	CB2-CA2-N2	-5.28	121.60	128.76
1	A	66	QFG	CB2-CA2-N2	-5.01	121.96	128.76
1	H	66	QFG	CB2-CA2-N2	-4.38	122.81	128.76
1	F	66	QFG	CB2-CA2-N2	-4.01	123.32	128.76
1	E	66	QFG	C2-CA2-N2	-3.87	106.18	108.95
1	C	66	QFG	CG1-CB1-CA1	-3.87	101.80	113.51
1	G	66	QFG	CG1-CB1-CA1	-3.61	102.56	113.51
1	A	66	QFG	CG1-CB1-CA1	-3.52	102.84	113.51
1	E	66	QFG	CG1-CB1-CA1	-3.33	103.41	113.51
1	C	66	QFG	CE2-CD2-CG2	-3.32	116.75	120.64
1	G	66	QFG	C3-CA3-N3	2.99	119.23	112.43
1	F	66	QFG	CG1-CB1-CA1	-2.97	104.52	113.51
1	H	66	QFG	O2-C2-N3	2.90	130.37	124.31
1	D	66	QFG	CA3-N3-C1	2.69	133.69	128.39
1	D	66	QFG	CG1-CB1-CA1	-2.53	105.85	113.51
1	D	66	QFG	CE1-CD1-CG2	-2.33	117.92	120.64
1	C	66	QFG	C2-CA2-N2	-2.29	107.31	108.95
1	B	66	QFG	CG1-CB1-CA1	-2.17	106.92	113.51
1	C	66	QFG	CZ-CE1-CD1	-2.17	117.56	120.24
1	C	66	QFG	CD2-CG2-CD1	2.10	120.76	117.65
1	D	66	QFG	CD2-CG2-CD1	2.10	120.76	117.65

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	66	QFG	C1-CA1-CB1-CG1

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Mol	Chain	Res	Type	Atoms
1	C	66	QFG	C1-CA1-CB1-CG1
1	E	66	QFG	C1-CA1-CB1-CG1
1	G	66	QFG	C1-CA1-CB1-CG1
1	A	66	QFG	C2-CA2-CB2-CG2
1	B	66	QFG	C2-CA2-CB2-CG2
1	C	66	QFG	C2-CA2-CB2-CG2
1	D	66	QFG	C2-CA2-CB2-CG2
1	E	66	QFG	C2-CA2-CB2-CG2
1	F	66	QFG	C2-CA2-CB2-CG2
1	G	66	QFG	C2-CA2-CB2-CG2
1	H	66	QFG	C2-CA2-CB2-CG2
1	A	66	QFG	N2-CA2-CB2-CG2
1	B	66	QFG	N2-CA2-CB2-CG2
1	C	66	QFG	N2-CA2-CB2-CG2
1	D	66	QFG	N2-CA2-CB2-CG2
1	E	66	QFG	N2-CA2-CB2-CG2
1	F	66	QFG	N2-CA2-CB2-CG2
1	G	66	QFG	N2-CA2-CB2-CG2
1	H	66	QFG	N2-CA2-CB2-CG2
1	A	66	QFG	C3-CA3-N3-C1
1	B	66	QFG	CA2-CB2-CG2-CD2
1	B	66	QFG	C1-CA1-CB1-CG1
1	D	66	QFG	C1-CA1-CB1-CG1
1	F	66	QFG	C1-CA1-CB1-CG1
1	H	66	QFG	C1-CA1-CB1-CG1
1	B	66	QFG	CA2-CB2-CG2-CD1
1	H	66	QFG	CA2-CB2-CG2-CD2
1	D	66	QFG	C3-CA3-N3-C1
1	E	66	QFG	C3-CA3-N3-C1
1	H	66	QFG	C3-CA3-N3-C1
1	F	66	QFG	CA2-CB2-CG2-CD1
1	H	66	QFG	CA2-CB2-CG2-CD1
1	D	66	QFG	CA2-CB2-CG2-CD1
1	C	66	QFG	CA2-CB2-CG2-CD1

There are no ring outliers.

8 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	66	QFG	3	0
1	E	66	QFG	1	0
1	B	66	QFG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	66	QFG	1	0
1	D	66	QFG	1	0
1	F	66	QFG	2	0
1	G	66	QFG	2	0
1	H	66	QFG	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 24 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	217/238 (91%)	-0.60	1 (0%) 87 85	12, 32, 44, 57	4 (1%)
1	B	218/238 (91%)	-0.61	3 (1%) 73 71	12, 30, 45, 56	3 (1%)
1	C	218/238 (91%)	-0.52	3 (1%) 73 71	13, 32, 46, 62	2 (0%)
1	D	218/238 (91%)	-0.60	2 (0%) 81 79	12, 29, 42, 54	3 (1%)
1	E	217/238 (91%)	-0.64	2 (0%) 81 79	13, 29, 42, 50	3 (1%)
1	F	218/238 (91%)	-0.60	2 (0%) 81 79	13, 30, 42, 53	2 (0%)
1	G	218/238 (91%)	-0.57	3 (1%) 73 71	20, 31, 44, 53	1 (0%)
1	H	217/238 (91%)	-0.60	2 (0%) 81 79	13, 33, 47, 63	2 (0%)
All	All	1741/1904 (91%)	-0.59	18 (1%) 79 77	12, 31, 45, 63	20 (1%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	225	ALA	5.0
1	F	6	SER	4.1
1	F	5	MET	4.0
1	B	225	ALA	4.0
1	H	225	ALA	3.8
1	C	5	MET	3.7
1	D	5	MET	3.5
1	B	5	MET	3.3
1	G	5	MET	3.2
1	B	6	SER	3.2
1	G	6	SER	3.2
1	C	225	ALA	3.1
1	A	225	ALA	2.9
1	D	225	ALA	2.7
1	C	6	SER	2.5
1	E	225	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	H	6	SER	2.2
1	E	6	SER	2.2

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	QFG	A	66	23/24	0.93	0.15	27,31,36,36	0
1	QFG	D	66	23/24	0.93	0.13	28,30,34,37	0
1	QFG	E	66	23/24	0.93	0.13	24,30,33,38	0
1	QFG	B	66	23/24	0.94	0.14	25,31,35,37	0
1	QFG	C	66	23/24	0.94	0.15	30,33,37,39	0
1	QFG	H	66	23/24	0.94	0.14	27,32,38,40	0
1	QFG	G	66	23/24	0.95	0.13	25,30,35,38	0
1	QFG	F	66	23/24	0.95	0.10	26,30,34,36	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [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
3	CL	B	228	1/1	0.96	0.22	37,37,37,37	0
3	CL	D	229	1/1	0.97	0.25	36,36,36,36	0
3	CL	F	226	1/1	0.97	0.13	26,26,26,26	0
3	CL	H	228	1/1	0.97	0.15	31,31,31,31	0
3	CL	F	227	1/1	0.98	0.16	34,34,34,34	0
3	CL	G	227	1/1	0.98	0.10	33,33,33,33	0
3	CL	H	227	1/1	0.98	0.16	31,31,31,31	0
3	CL	E	227	1/1	0.98	0.17	31,31,31,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CL	E	228	1/1	0.99	0.29	43,43,43,43	0
3	CL	C	228	1/1	0.99	0.19	32,32,32,32	0
3	CL	C	229	1/1	0.99	0.19	33,33,33,33	0
3	CL	G	226	1/1	0.99	0.18	33,33,33,33	0
3	CL	D	228	1/1	0.99	0.13	26,26,26,26	0
3	CL	B	227	1/1	0.99	0.14	28,28,28,28	0
3	CL	A	228	1/1	0.99	0.25	44,44,44,44	0
2	IOD	E	226	1/1	1.00	0.01	33,33,33,33	0
2	IOD	H	226	1/1	1.00	0.01	35,35,35,35	0
3	CL	A	227	1/1	1.00	0.15	29,29,29,29	0
2	IOD	A	226	1/1	1.00	0.02	33,33,33,33	0
2	IOD	B	226	1/1	1.00	0.01	35,35,35,35	0
2	IOD	C	226	1/1	1.00	0.01	34,34,34,34	0
2	IOD	C	227	1/1	1.00	0.01	33,33,33,33	0
2	IOD	D	226	1/1	1.00	0.01	32,32,32,32	0
2	IOD	D	227	1/1	1.00	0.01	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.