



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:03 PM UTC

PDB ID : 8C1X / pdb_00008c1x
Title : sfGFP C148 F206 mutant
Authors : Ahmed, R.D.; Rizkallah, P.J.; Jones, D.D.
Deposited on : 2022-12-21
Resolution : 1.89 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

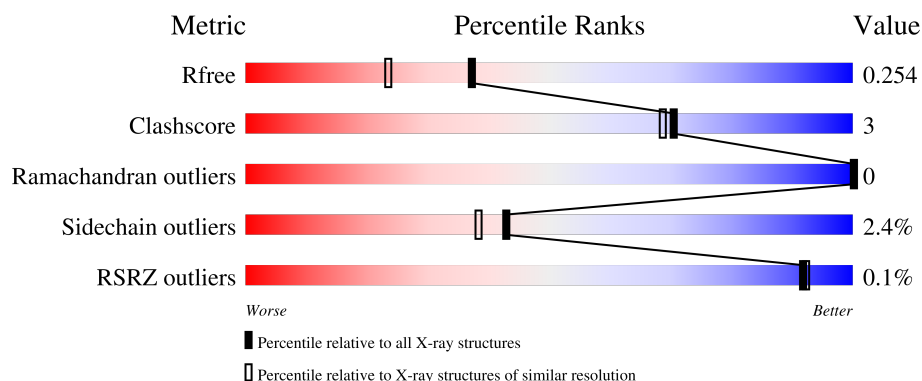
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	231	
1	B	231	
1	C	231	
1	D	231	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7679 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 Green fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	2	0
			1862	1177	321	356	8			
1	B	231	Total	C	N	O	S	0	3	0
			1870	1181	322	359	8			
1	C	229	Total	C	N	O	S	0	3	0
			1848	1170	317	353	8			
1	D	229	Total	C	N	O	S	0	3	0
			1850	1171	316	356	7			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ARG	SER	engineered mutation	UNP P42212
A	30	ARG	SER	engineered mutation	UNP P42212
A	39	ASN	TYR	engineered mutation	UNP P42212
A	64	LEU	PHE	engineered mutation	UNP P42212
A	66	CRO	SER	chromophore	UNP P42212
A	66	CRO	TYR	chromophore	UNP P42212
A	66	CRO	GLY	chromophore	UNP P42212
A	80	ARG	GLN	engineered mutation	UNP P42212
A	99	SER	PHE	engineered mutation	UNP P42212
A	105	THR	ASN	engineered mutation	UNP P42212
A	145	PHE	TYR	engineered mutation	UNP P42212
A	148	CYS	HIS	engineered mutation	UNP P42212
A	153	THR	MET	engineered mutation	UNP P42212
A	163	ALA	VAL	engineered mutation	UNP P42212
A	171	VAL	ILE	engineered mutation	UNP P42212
A	206	PHE	ALA	engineered mutation	UNP P42212
B	2	ARG	SER	engineered mutation	UNP P42212
B	30	ARG	SER	engineered mutation	UNP P42212
B	39	ASN	TYR	engineered mutation	UNP P42212
B	64	LEU	PHE	engineered mutation	UNP P42212
B	66	CRO	SER	chromophore	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
B	66	CRO	TYR	chromophore	UNP P42212
B	66	CRO	GLY	chromophore	UNP P42212
B	80	ARG	GLN	engineered mutation	UNP P42212
B	99	SER	PHE	engineered mutation	UNP P42212
B	105	THR	ASN	engineered mutation	UNP P42212
B	145	PHE	TYR	engineered mutation	UNP P42212
B	148	CYS	HIS	engineered mutation	UNP P42212
B	153	THR	MET	engineered mutation	UNP P42212
B	163	ALA	VAL	engineered mutation	UNP P42212
B	171	VAL	ILE	engineered mutation	UNP P42212
B	206	PHE	ALA	engineered mutation	UNP P42212
C	2	ARG	SER	engineered mutation	UNP P42212
C	30	ARG	SER	engineered mutation	UNP P42212
C	39	ASN	TYR	engineered mutation	UNP P42212
C	64	LEU	PHE	engineered mutation	UNP P42212
C	66	CRO	SER	chromophore	UNP P42212
C	66	CRO	TYR	chromophore	UNP P42212
C	66	CRO	GLY	chromophore	UNP P42212
C	80	ARG	GLN	engineered mutation	UNP P42212
C	99	SER	PHE	engineered mutation	UNP P42212
C	105	THR	ASN	engineered mutation	UNP P42212
C	145	PHE	TYR	engineered mutation	UNP P42212
C	148	CYS	HIS	engineered mutation	UNP P42212
C	153	THR	MET	engineered mutation	UNP P42212
C	163	ALA	VAL	engineered mutation	UNP P42212
C	171	VAL	ILE	engineered mutation	UNP P42212
C	206	PHE	ALA	engineered mutation	UNP P42212
D	2	ARG	SER	engineered mutation	UNP P42212
D	30	ARG	SER	engineered mutation	UNP P42212
D	39	ASN	TYR	engineered mutation	UNP P42212
D	64	LEU	PHE	engineered mutation	UNP P42212
D	66	CRO	SER	chromophore	UNP P42212
D	66	CRO	TYR	chromophore	UNP P42212
D	66	CRO	GLY	chromophore	UNP P42212
D	80	ARG	GLN	engineered mutation	UNP P42212
D	99	SER	PHE	engineered mutation	UNP P42212
D	105	THR	ASN	engineered mutation	UNP P42212
D	145	PHE	TYR	engineered mutation	UNP P42212
D	148	CYS	HIS	engineered mutation	UNP P42212
D	153	THR	MET	engineered mutation	UNP P42212
D	163	ALA	VAL	engineered mutation	UNP P42212
D	171	VAL	ILE	engineered mutation	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
D	206	PHE	ALA	engineered mutation	UNP P42212

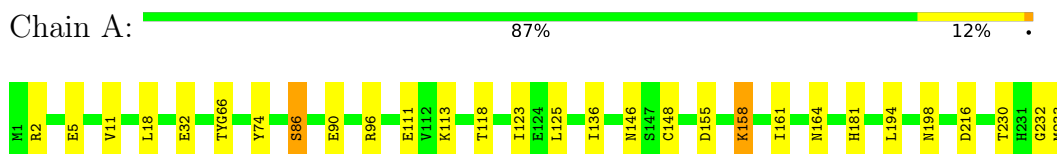
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	60	Total 60	O 60	0	0
2	B	62	Total 62	O 62	0	0
2	C	65	Total 65	O 65	0	0
2	D	62	Total 62	O 62	0	0

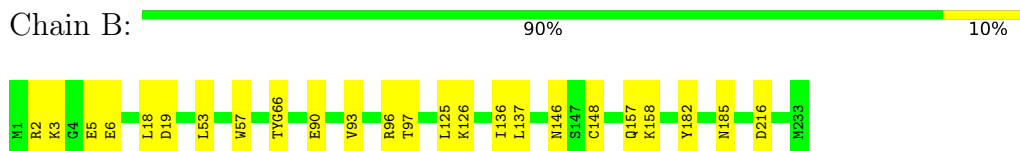
3 Residue-property plots [i](#)

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.

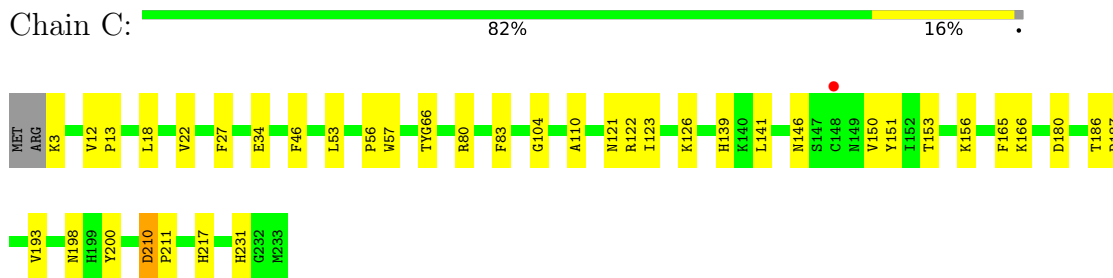
- Molecule 1: Green fluorescent protein



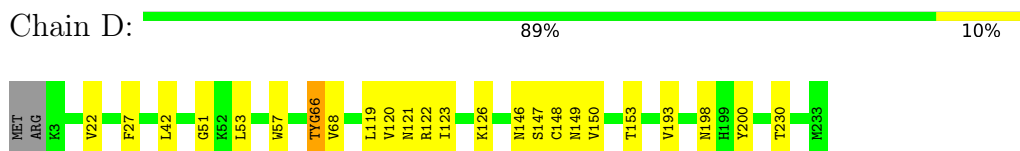
- Molecule 1: Green fluorescent protein



- Molecule 1: Green fluorescent protein



- Molecule 1: Green fluorescent protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	67.49Å 73.76Å 127.28Å 90.00° 90.11° 90.00°	Depositor
Resolution (Å)	39.24 – 1.89 39.24 – 1.89	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.24-1.89) 95.5 (39.24-1.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.00 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.206 , 0.247 0.212 , 0.254	Depositor DCC
R_{free} test set	4775 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	44.2	Xtriage
Anisotropy	0.381	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.470 for h,-k,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	7679	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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: CRO

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	1.14	2/1879 (0.1%)	1.40	7/2535 (0.3%)
1	B	1.13	0/1887	1.36	4/2546 (0.2%)
1	C	1.21	4/1865 (0.2%)	1.35	5/2518 (0.2%)
1	D	1.13	2/1867 (0.1%)	1.35	1/2521 (0.0%)
All	All	1.15	8/7498 (0.1%)	1.36	17/10120 (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	A	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	11	VAL	C-O	-6.14	1.17	1.24
1	C	139	HIS	CE1-NE2	6.11	1.38	1.32
1	C	104	GLY	C-O	5.93	1.29	1.23
1	D	51	GLY	C-O	5.89	1.27	1.24
1	D	193	VAL	C-O	-5.28	1.18	1.23
1	A	158	LYS	C-O	5.24	1.30	1.24
1	C	193	VAL	C-O	-5.22	1.17	1.23
1	C	187	PRO	C-O	-5.01	1.18	1.23

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	19	ASP	CA-C-N	7.17	127.13	122.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	19	ASP	C-N-CA	7.17	127.13	122.18
1	C	210	ASP	CB-CA-C	6.43	118.64	109.38
1	A	155	ASP	CA-CB-CG	6.04	118.64	112.60
1	A	118	THR	CA-CB-OG1	-6.00	100.60	109.60
1	C	186	THR	CB-CA-C	5.89	118.85	110.02
1	A	198	ASN	CA-CB-CG	-5.81	106.79	112.60
1	D	230	THR	CA-CB-OG1	-5.36	101.56	109.60
1	B	216	ASP	CA-CB-CG	5.21	117.81	112.60
1	A	216	ASP	CA-CB-CG	5.18	117.78	112.60
1	C	187	PRO	N-CA-CB	5.16	107.91	103.31
1	A	86	SER	N-CA-CB	5.14	118.14	110.22
1	A	74	TYR	CB-CA-C	5.13	116.77	109.38
1	A	230	THR	CA-CB-OG1	-5.09	101.97	109.60
1	B	97	THR	CA-CB-OG1	-5.02	102.07	109.60
1	C	34	GLU	CA-C-N	5.01	125.99	122.33
1	C	34	GLU	C-N-CA	5.01	125.99	122.33

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	96	ARG	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	1862	0	1814	13	0
1	B	1870	0	1817	10	0
1	C	1848	0	1797	16	0
1	D	1850	0	1794	11	0
2	A	60	0	0	1	0
2	B	62	0	0	1	0
2	C	65	0	0	2	0
2	D	62	0	0	0	0
All	All	7679	0	7222	50	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 (50) 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:2:ARG:NH1	1:A:86:SER:HB3	1.85	0.92
1:A:2:ARG:HH11	1:A:86:SER:HB3	1.37	0.88
1:A:2:ARG:NH1	1:A:194:LEU:HD12	2.02	0.74
1:B:2:ARG:HG2	1:B:5:GLU:HG2	1.70	0.73
1:D:120:VAL:HG11	1:D:122:ARG:NH1	2.04	0.71
1:C:151:TYR:HB2	2:C:302:HOH:O	1.96	0.63
1:D:66:CRO:OH	1:D:147:SER:HB3	2.06	0.54
1:A:18:LEU:C	1:A:18:LEU:HD23	2.32	0.54
1:A:2:ARG:NH1	1:A:86:SER:CB	2.67	0.53
1:A:232:GLY:O	1:A:233:MET:HB2	2.07	0.53
1:D:53:LEU:HD22	1:D:57:TRP:CE2	2.43	0.53
1:C:22:VAL:HG23	1:C:27:PHE:HE1	1.73	0.53
1:B:18:LEU:C	1:B:18:LEU:HD23	2.36	0.51
1:B:136:ILE:HD12	1:B:136:ILE:N	2.25	0.51
1:C:126:LYS:HE2	2:C:308:HOH:O	2.09	0.51
1:C:150:VAL:O	1:C:200:TYR:HA	2.13	0.49
1:C:150:VAL:HG13	1:C:165:PHE:CD1	2.47	0.48
1:D:22:VAL:HG23	1:D:27:PHE:HE1	1.78	0.48
1:C:53:LEU:HD22	1:C:57:TRP:CE2	2.48	0.48
1:B:93:VAL:O	1:B:185:ASN:HA	2.14	0.47
1:D:42[B]:LEU:HD21	1:D:68:VAL:HG23	1.96	0.47
1:B:125:LEU:C	1:B:125:LEU:HD23	2.41	0.45
1:C:56:PRO:HG2	1:C:141:LEU:HD12	1.98	0.45
1:C:18:LEU:C	1:C:18:LEU:HD23	2.42	0.45
1:D:150:VAL:O	1:D:200:TYR:HA	2.16	0.44
1:C:83:PHE:C	1:C:83:PHE:CD1	2.94	0.44
1:B:96:ARG:HA	1:B:182:TYR:O	2.18	0.44
1:A:161:ILE:C	1:A:161:ILE:HD12	2.43	0.44
1:C:110:ALA:HA	1:C:122:ARG:O	2.18	0.44
1:D:149:ASN:HB3	1:D:200:TYR:CD1	2.52	0.44
1:B:90:GLU:HG3	2:B:353:HOH:O	2.18	0.43
1:A:125:LEU:C	1:A:125:LEU:HD23	2.44	0.42
1:D:153:THR:HG22	1:D:198:ASN:OD1	2.19	0.42
1:C:46:PHE:O	1:C:217:HIS:HB2	2.19	0.42
1:A:136:ILE:HD12	1:A:136:ILE:N	2.35	0.42
1:A:32:GLU:OE2	2:A:301:HOH:O	2.22	0.42
1:D:119:LEU:HD13	1:D:119:LEU:C	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:VAL:HA	1:C:13:PRO:HD3	1.91	0.41
1:C:166:LYS:HG2	1:C:180:ASP:OD1	2.21	0.41
1:C:153:THR:HG22	1:C:198:ASN:OD1	2.21	0.41
1:D:121:ASN:ND2	1:D:123:ILE:HD11	2.36	0.41
1:A:111:GLU:OE1	1:A:113:LYS:NZ	2.54	0.41
1:A:164:ASN:HA	1:A:181:HIS:O	2.21	0.41
1:B:157:GLN:O	1:B:158:LYS:HB2	2.21	0.40
1:C:121:ASN:ND2	1:C:123:ILE:HD11	2.37	0.40
1:A:2:ARG:NH1	1:A:194:LEU:CD1	2.80	0.40
1:B:137:LEU:HD23	1:B:137:LEU:HA	1.87	0.40
1:D:53:LEU:HD22	1:D:57:TRP:CD2	2.55	0.40
1:B:53:LEU:HD22	1:B:57:TRP:CE2	2.56	0.40
1:C:210:ASP:HA	1:C:211:PRO:HD2	1.94	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	228/231 (99%)	219 (96%)	9 (4%)	0	100	100
1	B	229/231 (99%)	219 (96%)	10 (4%)	0	100	100
1	C	227/231 (98%)	222 (98%)	5 (2%)	0	100	100
1	D	227/231 (98%)	218 (96%)	9 (4%)	0	100	100
All	All	911/924 (99%)	878 (96%)	33 (4%)	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	203/201 (101%)	197 (97%)	6 (3%)	36	30
1	B	204/201 (102%)	199 (98%)	5 (2%)	42	37
1	C	202/201 (100%)	197 (98%)	5 (2%)	42	37
1	D	202/201 (100%)	199 (98%)	3 (2%)	57	56
All	All	811/804 (101%)	792 (98%)	19 (2%)	43	40

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLU
1	A	90	GLU
1	A	123	ILE
1	A	146	ASN
1	A	148	CYS
1	A	158	LYS
1	B	3	LYS
1	B	6	GLU
1	B	126	LYS
1	B	146	ASN
1	B	148	CYS
1	C	3	LYS
1	C	80	ARG
1	C	146	ASN
1	C	156	LYS
1	C	231	HIS
1	D	126	LYS
1	D	146	ASN
1	D	148	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	B	139	HIS
1	D	121	ASN
1	D	135	ASN
1	D	204	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	CRO	B	66	1	22,23,24	0.67	0	30,32,34	0.99	2 (6%)
1	CRO	A	66	1	22,23,24	0.69	0	30,32,34	1.06	2 (6%)
1	CRO	D	66	1	22,23,24	0.72	0	30,32,34	1.16	4 (13%)
1	CRO	C	66	1	22,23,24	0.70	0	30,32,34	1.18	4 (13%)

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	CRO	B	66	1	-	0/12/31/32	0/2/2/2
1	CRO	A	66	1	-	1/12/31/32	0/2/2/2
1	CRO	D	66	1	-	0/12/31/32	0/2/2/2
1	CRO	C	66	1	-	0/12/31/32	0/2/2/2

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CRO	O3-C3-CA3	-2.77	113.01	125.47
1	A	66	CRO	C2-N3-C1	2.55	109.25	108.07
1	C	66	CRO	OG1-CB1-CA1	2.38	114.08	109.00
1	B	66	CRO	O3-C3-CA3	-2.36	114.84	125.47
1	C	66	CRO	CG2-CB2-CA2	2.20	132.48	129.87
1	C	66	CRO	O3-C3-CA3	-2.20	115.59	125.47
1	D	66	CRO	CA2-C2-N3	2.18	105.33	103.50
1	D	66	CRO	OG1-CB1-CA1	2.14	113.55	109.00
1	D	66	CRO	C3-CA3-N3	2.08	117.16	112.43
1	B	66	CRO	C2-CA2-N2	-2.06	107.48	108.95
1	D	66	CRO	C2-CA2-N2	-2.06	107.48	108.95
1	C	66	CRO	C1-CA1-N1	2.03	113.41	109.78

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	66	CRO	N3-C1-CA1-CB1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	66	CRO	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

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	230/231 (99%)	-1.07	0 100 100	22, 52, 81, 98	2 (0%)
1	B	230/231 (99%)	-1.06	0 100 100	21, 52, 85, 100	3 (1%)
1	C	228/231 (98%)	-1.06	1 (0%) 88 90	22, 50, 74, 92	3 (1%)
1	D	228/231 (98%)	-1.07	0 100 100	22, 51, 76, 101	3 (1%)
All	All	916/924 (99%)	-1.07	1 (0%) 92 92	21, 51, 80, 101	11 (1%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	148[A]	CYS	2.4

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	CRO	A	66	22/23	0.99	0.03	38,41,46,47	0
1	CRO	B	66	22/23	0.99	0.03	37,41,46,46	0
1	CRO	D	66	22/23	0.99	0.03	38,40,44,48	0
1	CRO	C	66	22/23	1.00	0.03	36,40,45,49	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.