



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 5, 2026 – 12:56 PM UTC

PDB ID : 3OGO / pdb_00003ogo
Title : Structure of the GFP:GFP-nanobody complex at 2.8 Å resolution in space-group P21212
Authors : Kubala, M.H.; Kovtun, O.; Alexandrov, K.; Collins, B.M.
Deposited on : 2010-08-17
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

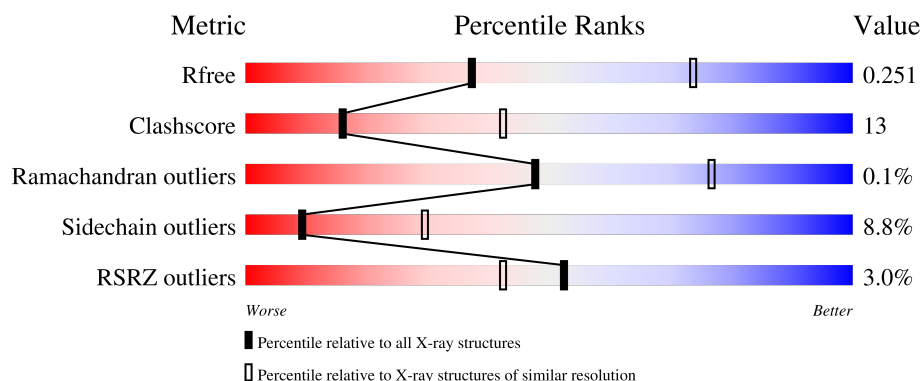
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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.



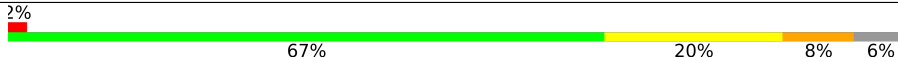
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	247	2% 74% 18% • 6%
1	B	247	2% 72% 17% • 9%
1	C	247	3% 62% 26% • 8%
1	D	247	3% 65% 23% • 9%
2	E	123	4% 65% 22% 7% 7%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	F	123	
2	G	123	
2	H	123	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	CRO	A	65	-	-	X	-
1	CRO	B	65	-	-	X	-
1	CRO	C	65	-	-	X	-
1	CRO	D	65	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11212 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	232	Total	C	N	O	S	0	0	0
			1853	1177	311	358	7			
1	B	225	Total	C	N	O	S	0	0	0
			1803	1148	304	345	6			
1	C	228	Total	C	N	O	S	0	1	0
			1833	1169	307	351	6			
1	D	226	Total	C	N	O	S	0	0	0
			1810	1152	305	347	6			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	ALA	-	expression tag	UNP P42212
A	-8	HIS	-	expression tag	UNP P42212
A	-7	HIS	-	expression tag	UNP P42212
A	-6	HIS	-	expression tag	UNP P42212
A	-5	HIS	-	expression tag	UNP P42212
A	-4	HIS	-	expression tag	UNP P42212
A	-3	HIS	-	expression tag	UNP P42212
A	-2	SER	-	expression tag	UNP P42212
A	-1	SER	-	expression tag	UNP P42212
A	0	GLY	-	expression tag	UNP P42212
A	1	VAL	-	expression tag	UNP P42212
A	64	LEU	PHE	engineered mutation	UNP P42212
A	65	CRO	SER	chromophore	UNP P42212
A	65	CRO	TYR	chromophore	UNP P42212
A	65	CRO	GLY	chromophore	UNP P42212
A	231	LEU	HIS	engineered mutation	UNP P42212
B	-9	ALA	-	expression tag	UNP P42212
B	-8	HIS	-	expression tag	UNP P42212
B	-7	HIS	-	expression tag	UNP P42212
B	-6	HIS	-	expression tag	UNP P42212
B	-5	HIS	-	expression tag	UNP P42212

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	HIS	-	expression tag	UNP P42212
B	-3	HIS	-	expression tag	UNP P42212
B	-2	SER	-	expression tag	UNP P42212
B	-1	SER	-	expression tag	UNP P42212
B	0	GLY	-	expression tag	UNP P42212
B	1	VAL	-	expression tag	UNP P42212
B	64	LEU	PHE	engineered mutation	UNP P42212
B	65	CRO	SER	chromophore	UNP P42212
B	65	CRO	TYR	chromophore	UNP P42212
B	65	CRO	GLY	chromophore	UNP P42212
B	231	LEU	HIS	engineered mutation	UNP P42212
C	-9	ALA	-	expression tag	UNP P42212
C	-8	HIS	-	expression tag	UNP P42212
C	-7	HIS	-	expression tag	UNP P42212
C	-6	HIS	-	expression tag	UNP P42212
C	-5	HIS	-	expression tag	UNP P42212
C	-4	HIS	-	expression tag	UNP P42212
C	-3	HIS	-	expression tag	UNP P42212
C	-2	SER	-	expression tag	UNP P42212
C	-1	SER	-	expression tag	UNP P42212
C	0	GLY	-	expression tag	UNP P42212
C	1	VAL	-	expression tag	UNP P42212
C	64	LEU	PHE	engineered mutation	UNP P42212
C	65	CRO	SER	chromophore	UNP P42212
C	65	CRO	TYR	chromophore	UNP P42212
C	65	CRO	GLY	chromophore	UNP P42212
C	231	LEU	HIS	engineered mutation	UNP P42212
D	-9	ALA	-	expression tag	UNP P42212
D	-8	HIS	-	expression tag	UNP P42212
D	-7	HIS	-	expression tag	UNP P42212
D	-6	HIS	-	expression tag	UNP P42212
D	-5	HIS	-	expression tag	UNP P42212
D	-4	HIS	-	expression tag	UNP P42212
D	-3	HIS	-	expression tag	UNP P42212
D	-2	SER	-	expression tag	UNP P42212
D	-1	SER	-	expression tag	UNP P42212
D	0	GLY	-	expression tag	UNP P42212
D	1	VAL	-	expression tag	UNP P42212
D	64	LEU	PHE	engineered mutation	UNP P42212
D	65	CRO	SER	chromophore	UNP P42212
D	65	CRO	TYR	chromophore	UNP P42212
D	65	CRO	GLY	chromophore	UNP P42212

Continued on next page...

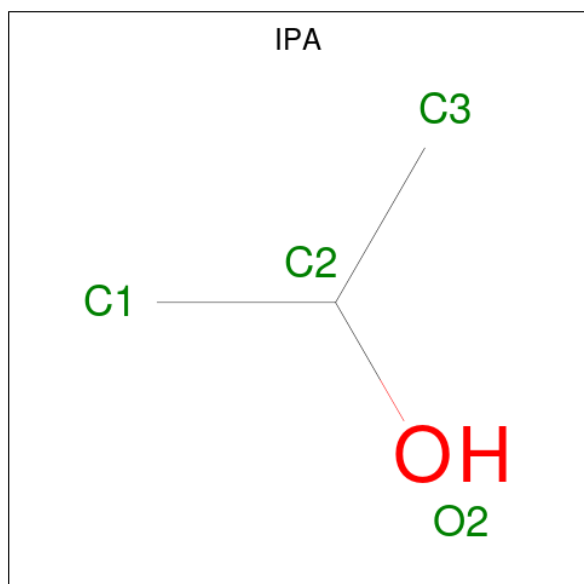
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	231	LEU	HIS	engineered mutation	UNP P42212

- Molecule 2 is a protein called GFP-nanobody.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	115	Total	C	N	O	S	0	0	0
			893	553	160	175	5			
2	F	115	Total	C	N	O	S	0	0	0
			893	553	160	175	5			
2	G	116	Total	C	N	O	S	0	0	0
			901	558	161	176	6			
2	H	116	Total	C	N	O	S	0	0	0
			902	559	162	176	5			

- Molecule 3 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: C_3H_8O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	3	1		
3	D	1	Total	C	O	0	0
			4	3	1		
3	F	1	Total	C	O	0	0
			4	3	1		

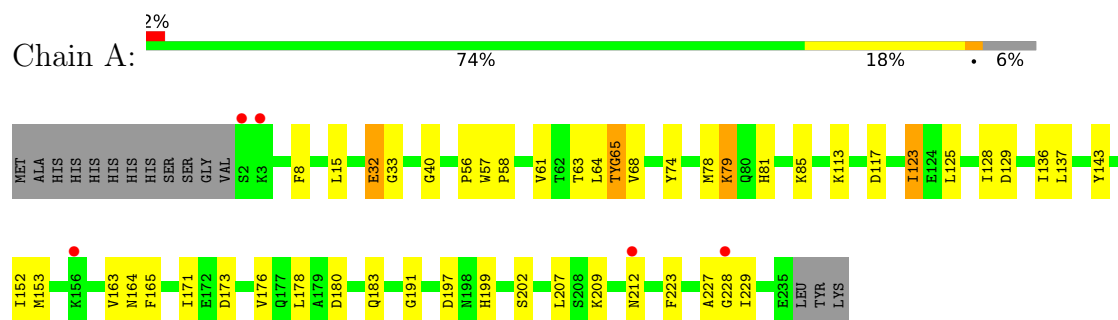
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	65	Total 65	O 65	0	0
4	B	63	Total 63	O 63	0	0
4	C	45	Total 45	O 45	0	0
4	D	42	Total 42	O 42	0	0
4	E	22	Total 22	O 22	0	0
4	F	25	Total 25	O 25	0	0
4	G	29	Total 29	O 29	0	0
4	H	21	Total 21	O 21	0	0

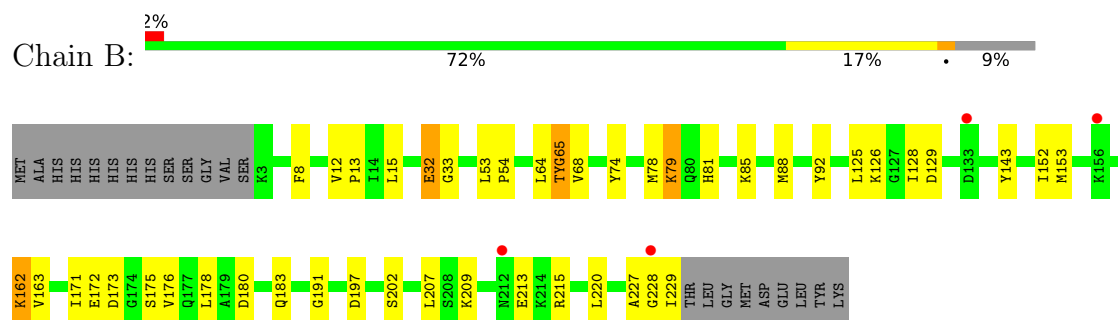
3 Residue-property plots [i](#)

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.

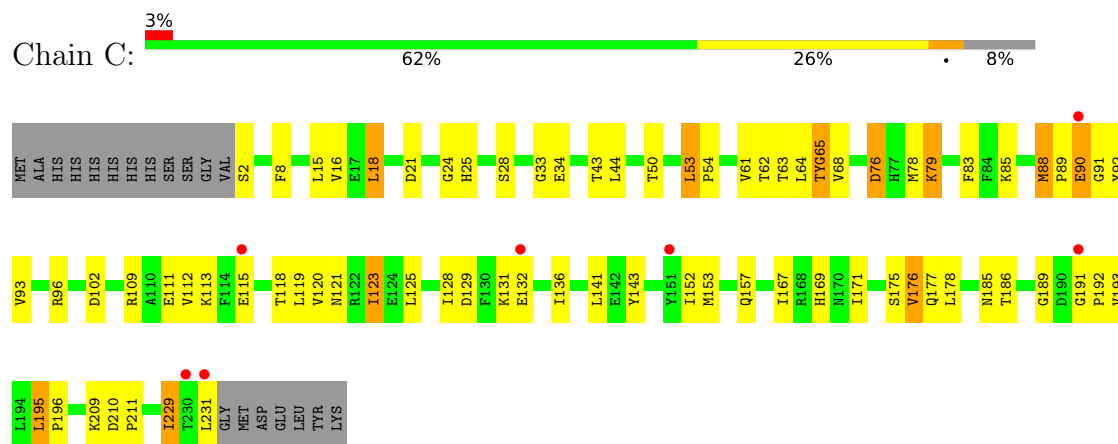
- Molecule 1: Green fluorescent protein

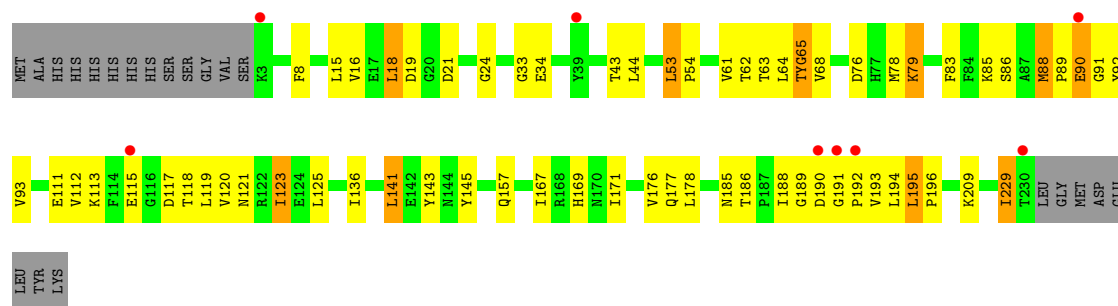


- Molecule 1: Green fluorescent protein



- Molecule 1: Green fluorescent protein

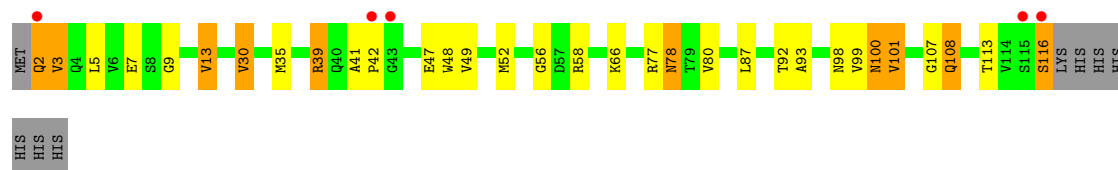




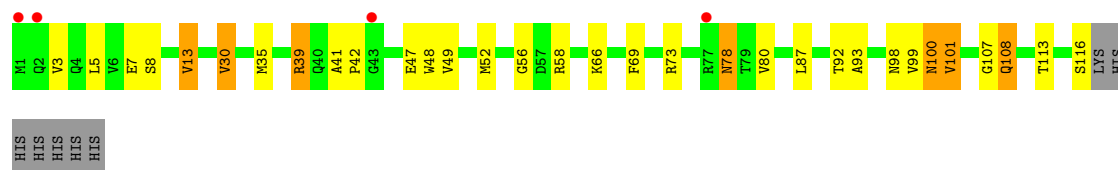
• Molecule 2: GFP-nanobody



• Molecule 2: GFP-nanobody



• Molecule 2: GFP-nanobody



• Molecule 2: GFP-nanobody



HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	140.10Å 147.60Å 101.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.43 – 2.80 42.43 – 2.80	Depositor EDS
% Data completeness (in resolution range)	87.0 (42.43-2.80) 87.0 (42.43-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.02 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.202 , 0.253 0.199 , 0.251	Depositor DCC
R_{free} test set	2433 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	37.2	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11212	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 4.23% 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: IPA, 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	0.54	0/1871	0.83	0/2527
1	B	0.56	0/1821	0.85	0/2460
1	C	0.49	0/1855	0.85	4/2507 (0.2%)
1	D	0.51	0/1828	0.85	4/2470 (0.2%)
2	E	0.49	0/912	0.79	0/1234
2	F	0.51	0/912	0.80	1/1234 (0.1%)
2	G	0.49	0/920	0.80	0/1244
2	H	0.52	0/921	0.79	0/1245
All	All	0.52	0/11040	0.83	9/14921 (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

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	195	LEU	CA-C-N	6.20	126.11	119.85
1	D	195	LEU	C-N-CA	6.20	126.11	119.85
1	C	195	LEU	CA-C-N	6.05	125.96	119.85
1	C	195	LEU	C-N-CA	6.05	125.96	119.85
2	F	3	VAL	N-CA-C	6.02	117.39	109.21

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	228	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	1853	0	1799	39	0
1	B	1803	0	1754	36	0
1	C	1833	0	1786	68	0
1	D	1810	0	1761	53	0
2	E	893	0	848	27	0
2	F	893	0	848	25	0
2	G	901	0	860	22	0
2	H	902	0	861	25	0
3	A	4	0	8	0	0
3	D	4	0	8	1	0
3	F	4	0	8	0	0
4	A	65	0	0	2	0
4	B	63	0	0	1	0
4	C	45	0	0	4	0
4	D	42	0	0	4	0
4	E	22	0	0	2	0
4	F	25	0	0	0	0
4	G	29	0	0	2	0
4	H	21	0	0	1	0
All	All	11212	0	10541	283	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 13.

The worst 5 of 283 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:CRO:C3	1:B:68:VAL:N	2.00	1.23
1:B:64:LEU:C	1:B:65:CRO:H2	1.55	1.12
1:B:64:LEU:O	1:B:65:CRO:N1	1.94	0.99
3:D:239:IPA:H32	2:F:9:GLY:H	1.33	0.93

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:CRO:C3	1:A:68:VAL:N	2.32	0.93

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	227/247 (92%)	217 (96%)	9 (4%)	1 (0%)	30	60
1	B	220/247 (89%)	209 (95%)	10 (4%)	1 (0%)	24	55
1	C	224/247 (91%)	218 (97%)	6 (3%)	0	100	100
1	D	221/247 (90%)	211 (96%)	10 (4%)	0	100	100
2	E	113/123 (92%)	110 (97%)	3 (3%)	0	100	100
2	F	113/123 (92%)	110 (97%)	3 (3%)	0	100	100
2	G	114/123 (93%)	111 (97%)	3 (3%)	0	100	100
2	H	114/123 (93%)	111 (97%)	3 (3%)	0	100	100
All	All	1346/1480 (91%)	1297 (96%)	47 (4%)	2 (0%)	48	77

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	191	GLY
1	B	191	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	202/215 (94%)	191 (95%)	11 (5%)	20	51
1	B	196/215 (91%)	186 (95%)	10 (5%)	21	54
1	C	200/215 (93%)	184 (92%)	16 (8%)	11	34
1	D	197/215 (92%)	183 (93%)	14 (7%)	13	39
2	E	95/103 (92%)	81 (85%)	14 (15%)	3	10
2	F	95/103 (92%)	82 (86%)	13 (14%)	3	13
2	G	96/103 (93%)	83 (86%)	13 (14%)	4	13
2	H	96/103 (93%)	83 (86%)	13 (14%)	4	13
All	All	1177/1272 (92%)	1073 (91%)	104 (9%)	9	29

5 of 104 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	E	58	ARG
2	F	66	LYS
2	H	98	ASN
2	E	78	ASN
2	F	2	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 34 such sidechains are listed below:

Mol	Chain	Res	Type
2	H	4	GLN
2	H	78	ASN
2	H	108	GLN
2	E	111	GLN
2	E	108	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	65	-	22,23,24	3.10	9 (40%)	30,32,34	2.61	9 (30%)
1	CRO	A	65	-	22,23,24	3.13	10 (45%)	30,32,34	2.52	8 (26%)
1	CRO	D	65	-	22,23,24	3.19	10 (45%)	30,32,34	2.24	9 (30%)
1	CRO	C	65	-	22,23,24	3.01	10 (45%)	30,32,34	2.54	9 (30%)

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	65	-	-	0/12/31/32	0/2/2/2
1	CRO	A	65	-	-	2/12/31/32	0/2/2/2
1	CRO	D	65	-	-	0/12/31/32	0/2/2/2
1	CRO	C	65	-	-	2/12/31/32	0/2/2/2

The worst 5 of 39 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	65	CRO	C1-N2	9.28	1.45	1.32
1	B	65	CRO	C1-N2	8.87	1.45	1.32
1	D	65	CRO	C1-N2	8.79	1.45	1.32
1	C	65	CRO	C1-N2	8.45	1.44	1.32
1	B	65	CRO	C1-N3	6.56	1.48	1.37

The worst 5 of 35 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	65	CRO	O2-C2-CA2	-8.36	125.68	131.02
1	B	65	CRO	O2-C2-CA2	-7.86	126.01	131.02
1	A	65	CRO	O2-C2-CA2	-7.40	126.30	131.02
1	D	65	CRO	O2-C2-CA2	-5.98	127.21	131.02
1	C	65	CRO	C2-N3-C1	-5.50	105.52	108.07

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	65	CRO	C3-CA3-N3-C2
1	C	65	CRO	C3-CA3-N3-C2
1	A	65	CRO	C3-CA3-N3-C1
1	C	65	CRO	C3-CA3-N3-C1

There are no ring outliers.

4 monomers are involved in 53 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	65	CRO	10	0
1	A	65	CRO	14	0
1	D	65	CRO	12	0
1	C	65	CRO	17	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands 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	$\# Z > 2$
3	IPA	D	239	-	3,3,3	0.60	0	3,3,3	0.21	0
3	IPA	F	124	-	3,3,3	0.69	0	3,3,3	0.36	0
3	IPA	A	239	-	3,3,3	0.67	0	3,3,3	0.09	0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	239	IPA	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	2
1	A	2
1	D	2
1	B	2

The worst 5 of 8 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	65:CRO	C3	68:VAL	N	2.65
1	A	64:LEU	C	65:CRO	N1	2.40
1	D	65:CRO	C3	68:VAL	N	2.38
1	D	64:LEU	C	65:CRO	N1	2.36
1	A	65:CRO	C3	68:VAL	N	2.32

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	231/247 (93%)	-0.42	5 (2%) 62 52	15, 26, 56, 88	0
1	B	224/247 (90%)	-0.44	4 (1%) 67 58	15, 26, 54, 83	0
1	C	227/247 (91%)	-0.04	7 (3%) 51 41	20, 37, 68, 82	1 (0%)
1	D	225/247 (91%)	0.12	8 (3%) 46 37	20, 39, 63, 89	0
2	E	115/123 (93%)	-0.22	5 (4%) 40 31	19, 31, 60, 77	0
2	F	115/123 (93%)	0.00	5 (4%) 40 31	21, 31, 61, 82	0
2	G	116/123 (94%)	-0.15	4 (3%) 48 39	22, 30, 63, 76	0
2	H	116/123 (94%)	-0.11	3 (2%) 57 47	21, 30, 62, 77	0
All	All	1369/1480 (92%)	-0.17	41 (2%) 52 42	15, 32, 63, 89	1 (0%)

The worst 5 of 41 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	43	GLY	5.3
1	D	191	GLY	4.3
2	G	2	GLN	4.2
2	F	42	PRO	4.2
1	D	192	PRO	3.8

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

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	D	65	22/23	0.94	0.09	18,32,43,49	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	CRO	C	65	22/23	0.96	0.08	19,28,37,38	0
1	CRO	B	65	22/23	0.96	0.07	16,21,27,40	0
1	CRO	A	65	22/23	0.97	0.06	10,20,27,29	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($\text{\AA}^2$)	Q<0.9
3	IPA	F	124	4/4	0.55	0.27	17,39,48,71	0
3	IPA	A	239	4/4	0.76	0.25	32,42,49,61	0
3	IPA	D	239	4/4	0.86	0.17	31,32,33,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.