



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 01:59 PM UTC

PDB ID : 1F0B / pdb_00001f0b
Title : CRYSTAL STRUCTURE OF THE GREEN FLUORESCENT PROTEIN
(GFP) VARIANT YFP-H148Q
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Deposited on : 2000-05-15
Resolution : 2.10 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

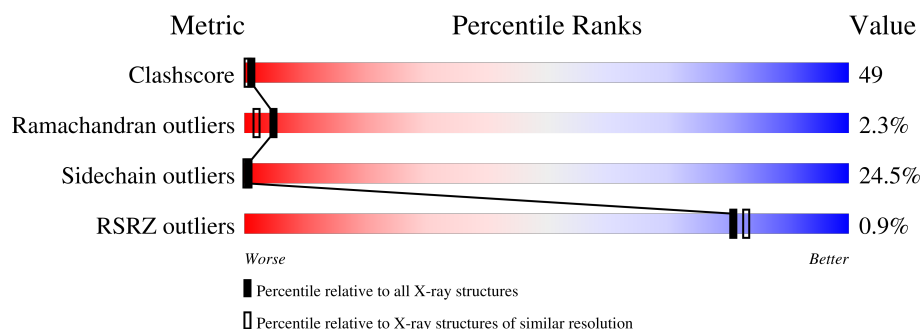
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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))
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	236	29% 38% 24% • •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1941 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	226	Total	C	N	O	S	36	0	0
			1816	1159	306	345	6			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	66	CR2	SER	chromophore	UNP P42212
A	66	CR2	TYR	chromophore	UNP P42212
A	66	CR2	GLY	chromophore	UNP P42212
A	68	LEU	VAL	engineered mutation	UNP P42212
A	72	ALA	SER	engineered mutation	UNP P42212
A	80	ARG	GLN	engineered mutation	UNP P42212
A	148	GLN	HIS	engineered mutation	UNP P42212
A	203	TYR	THR	engineered mutation	UNP P42212

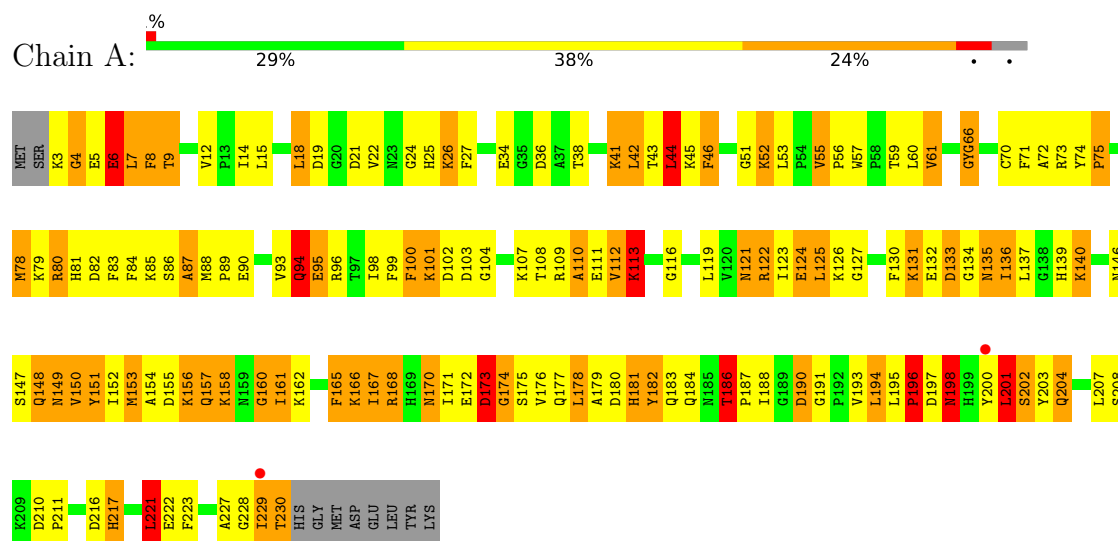
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	125	Total	O	0	0
			125	125		

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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.66Å 62.65Å 66.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.50 – 2.10 22.50 – 2.10	Depositor EDS
% Data completeness (in resolution range)	81.5 (22.50-2.10) 82.2 (22.50-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.82 (at 2.04Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.188 , (Not available) 0.197 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 137.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	1941	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.61% 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: CR2

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	1.69	14/1838 (0.8%)	2.08	63/2482 (2.5%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	139	HIS	CA-C	7.27	1.62	1.52
1	A	124	GLU	CA-C	-6.83	1.44	1.52
1	A	46	PHE	CA-C	-6.44	1.44	1.52
1	A	99	PHE	N-CA	-6.21	1.38	1.46
1	A	78	MET	CA-C	-5.92	1.45	1.52
1	A	184	GLN	CA-C	-5.76	1.45	1.52
1	A	57	TRP	N-CA	5.72	1.52	1.46
1	A	110	ALA	CA-C	-5.50	1.45	1.52
1	A	55	VAL	C-N	-5.41	1.27	1.33
1	A	93	VAL	C-N	-5.36	1.26	1.33
1	A	87	ALA	C-O	-5.29	1.17	1.24
1	A	8	PHE	C-N	-5.24	1.27	1.33
1	A	221	LEU	CA-C	-5.08	1.46	1.52
1	A	121	ASN	CA-C	-5.00	1.47	1.52

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	ASP	CA-CB-CG	11.91	124.51	112.60
1	A	191	GLY	N-CA-C	10.73	124.88	112.00
1	A	100	PHE	CA-CB-CG	-9.55	104.25	113.80
1	A	157	GLN	N-CA-C	-8.81	100.06	112.13
1	A	46	PHE	CA-CB-CG	-8.78	105.02	113.80
1	A	135	ASN	CA-CB-CG	-8.15	104.44	112.60
1	A	41	LYS	N-CA-C	7.95	122.65	109.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	HIS	CA-CB-CG	-7.85	105.95	113.80
1	A	149	ASN	CA-CB-CG	-7.80	104.80	112.60
1	A	165	PHE	CA-CB-CG	-7.66	106.14	113.80
1	A	201	LEU	N-CA-CB	7.65	122.89	110.42
1	A	150	VAL	N-CA-CB	-7.36	101.81	111.82
1	A	9	THR	N-CA-CB	-7.23	99.21	110.06
1	A	61	VAL	N-CA-C	7.11	117.24	110.42
1	A	102	ASP	CA-CB-CG	6.89	119.49	112.60
1	A	198	ASN	CA-CB-CG	-6.85	105.75	112.60
1	A	72	ALA	CA-C-N	-6.78	114.02	122.84
1	A	72	ALA	C-N-CA	-6.78	114.02	122.84
1	A	36	ASP	CA-C-N	6.78	129.67	120.38
1	A	36	ASP	C-N-CA	6.78	129.67	120.38
1	A	36	ASP	CA-CB-CG	6.76	119.36	112.60
1	A	84	PHE	N-CA-C	6.71	118.25	111.07
1	A	26	LYS	N-CA-C	6.70	119.61	108.96
1	A	78	MET	CB-CA-C	-6.68	100.27	111.02
1	A	75	PRO	N-CA-CB	6.65	109.13	103.35
1	A	116	GLY	N-CA-C	-6.62	100.08	111.47
1	A	122	ARG	N-CA-CB	-6.61	100.19	110.57
1	A	197	ASP	N-CA-C	-6.51	101.76	110.55
1	A	94	GLN	N-CA-CB	-6.41	100.50	110.57
1	A	191	GLY	CA-C-O	6.34	128.28	121.62
1	A	12	VAL	CA-C-N	6.16	126.67	120.14
1	A	12	VAL	C-N-CA	6.16	126.67	120.14
1	A	148	GLN	N-CA-C	6.07	117.77	108.42
1	A	173	ASP	CA-C-N	6.03	133.24	121.41
1	A	173	ASP	C-N-CA	6.03	133.24	121.41
1	A	201	LEU	CB-CA-C	6.00	119.92	110.19
1	A	26	LYS	CB-CA-C	6.00	120.24	110.22
1	A	196	PRO	CB-CA-C	-5.95	101.54	110.17
1	A	186	THR	N-CA-CB	5.91	121.21	111.17
1	A	208	SER	CA-C-N	-5.81	114.22	122.94
1	A	208	SER	C-N-CA	-5.81	114.22	122.94
1	A	14	ILE	CB-CA-C	-5.68	100.89	110.71
1	A	99	PHE	CB-CA-C	-5.64	101.06	110.19
1	A	182	TYR	N-CA-CB	5.55	119.27	110.84
1	A	156	LYS	N-CA-C	5.54	119.58	113.21
1	A	207	LEU	N-CA-CB	5.51	119.61	110.63
1	A	160	GLY	N-CA-C	-5.49	100.16	113.18
1	A	71	PHE	CA-CB-CG	-5.46	108.34	113.80
1	A	112	VAL	N-CA-C	5.46	115.96	107.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	LEU	CB-CA-C	-5.46	98.07	110.07
1	A	42	LEU	CB-CA-C	-5.41	99.80	109.71
1	A	45	LYS	N-CA-CB	5.39	119.86	110.81
1	A	8	PHE	CA-C-O	5.38	125.94	119.59
1	A	217	HIS	CA-CB-CG	-5.38	108.42	113.80
1	A	190	ASP	CA-C-N	5.26	128.09	121.36
1	A	190	ASP	C-N-CA	5.26	128.09	121.36
1	A	113	LYS	N-CA-C	5.18	116.39	108.42
1	A	6	GLU	N-CA-C	5.11	117.59	111.71
1	A	167	ILE	O-C-N	-5.05	117.99	123.14
1	A	173	ASP	N-CA-C	-5.03	99.43	108.58
1	A	27	PHE	CA-CB-CG	-5.02	108.78	113.80
1	A	101	LYS	CA-C-N	-5.00	115.72	123.17
1	A	101	LYS	C-N-CA	-5.00	115.72	123.17

There are no chirality outliers.

There are no planarity outliers.

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	1816	0	1762	171	0
2	A	125	0	0	12	0
All	All	1941	0	1762	171	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 49.

All (171) 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:198:ASN:H	1:A:198:ASN:HD22	1.03	0.97
1:A:198:ASN:H	1:A:198:ASN:ND2	1.59	0.96
1:A:152:ILE:HD12	1:A:200:TYR:HA	1.51	0.91
1:A:7:LEU:HD12	1:A:7:LEU:H	1.33	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:ASN:HA	1:A:140:LYS:HG2	1.55	0.88
1:A:166:LYS:HG3	1:A:178:LEU:HD21	1.56	0.88
1:A:146:ASN:HD21	1:A:170:ASN:HD21	1.19	0.86
1:A:135:ASN:HA	1:A:140:LYS:CG	2.05	0.86
1:A:108:THR:HG22	1:A:125:LEU:HB2	1.60	0.84
1:A:168:ARG:HB3	1:A:176:VAL:HG11	1.59	0.84
1:A:41:LYS:HG3	1:A:223:PHE:CE1	2.13	0.83
1:A:155:ASP:HB2	1:A:162:LYS:HG3	1.62	0.80
1:A:198:ASN:ND2	1:A:198:ASN:N	2.30	0.79
1:A:24:GLY:HA2	2:A:373:HOH:O	1.83	0.79
1:A:103:ASP:HA	1:A:131:LYS:HZ2	1.48	0.78
1:A:103:ASP:HA	1:A:131:LYS:NZ	2.01	0.76
1:A:22:VAL:HG22	1:A:127:GLY:HA3	1.67	0.75
1:A:7:LEU:HD12	1:A:7:LEU:N	2.01	0.74
1:A:155:ASP:CG	1:A:162:LYS:HE2	2.14	0.73
1:A:221:LEU:HD11	1:A:223:PHE:CE2	2.23	0.72
1:A:75:PRO:HA	2:A:431:HOH:O	1.90	0.71
1:A:7:LEU:H	1:A:7:LEU:CD1	2.04	0.71
1:A:119:LEU:HD12	2:A:433:HOH:O	1.92	0.70
1:A:78:MET:HG2	1:A:229:ILE:HG21	1.73	0.70
1:A:86:SER:O	1:A:193:VAL:HG12	1.94	0.68
1:A:152:ILE:CD1	1:A:200:TYR:HA	2.23	0.68
1:A:166:LYS:HD3	1:A:178:LEU:HD11	1.74	0.68
1:A:146:ASN:ND2	1:A:168:ARG:HH12	1.93	0.66
1:A:228:GLY:HA2	2:A:419:HOH:O	1.94	0.66
1:A:52:LYS:HE2	1:A:216:ASP:HB2	1.77	0.65
1:A:195:LEU:N	1:A:195:LEU:HD22	2.12	0.65
1:A:110:ALA:HB1	1:A:121:ASN:OD1	1.96	0.65
1:A:78:MET:HG2	1:A:229:ILE:CG2	2.26	0.65
1:A:135:ASN:HA	1:A:140:LYS:HG3	1.78	0.65
1:A:4:GLY:O	1:A:7:LEU:HD13	1.97	0.64
1:A:146:ASN:ND2	1:A:168:ARG:NH1	2.45	0.64
1:A:94:GLN:HG3	2:A:252:HOH:O	1.96	0.64
1:A:153:MET:HB2	1:A:198:ASN:OD1	1.98	0.64
1:A:146:ASN:ND2	1:A:170:ASN:HD21	1.94	0.64
1:A:104:GLY:HA2	2:A:284:HOH:O	1.97	0.63
1:A:146:ASN:HD22	1:A:168:ARG:NH1	1.96	0.63
1:A:55:VAL:HB	1:A:56:PRO:HD2	1.81	0.63
1:A:166:LYS:CD	1:A:178:LEU:HD11	2.29	0.63
1:A:155:ASP:OD2	1:A:158:LYS:HB2	1.98	0.63
1:A:125:LEU:HD12	1:A:126:LYS:N	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ASN:HD21	1:A:170:ASN:ND2	1.94	0.62
1:A:178:LEU:HD22	1:A:179:ALA:N	2.15	0.62
1:A:78:MET:CE	1:A:229:ILE:H	2.13	0.61
1:A:178:LEU:HD22	1:A:179:ALA:H	1.65	0.61
1:A:173:ASP:C	1:A:175:SER:H	2.03	0.61
1:A:148:GLN:HG3	1:A:149:ASN:N	2.16	0.61
1:A:70:CYS:O	1:A:85:LYS:HE3	2.01	0.60
1:A:151:TYR:CD2	1:A:200:TYR:CD1	2.89	0.60
1:A:21:ASP:OD2	1:A:26:LYS:HD3	2.02	0.60
1:A:168:ARG:HB3	1:A:176:VAL:CG1	2.30	0.59
1:A:21:ASP:HA	1:A:25:HIS:O	2.02	0.59
1:A:18:LEU:HD13	1:A:19:ASP:C	2.28	0.59
1:A:89:PRO:HD2	1:A:90:GLU:OE2	2.02	0.59
1:A:171:ILE:HG22	1:A:173:ASP:H	1.68	0.59
1:A:130:PHE:HE2	1:A:136:ILE:HD13	1.67	0.59
1:A:204:GLN:HG3	2:A:365:HOH:O	2.03	0.59
1:A:152:ILE:HD11	1:A:201:LEU:HG	1.85	0.58
1:A:100:PHE:CG	1:A:136:ILE:HD11	2.37	0.58
1:A:151:TYR:N	1:A:151:TYR:CD1	2.72	0.57
1:A:52:LYS:HE3	1:A:52:LYS:N	2.19	0.57
1:A:4:GLY:HA2	1:A:7:LEU:HD13	1.85	0.57
1:A:43:THR:HB	2:A:400:HOH:O	2.05	0.57
1:A:46:PHE:O	1:A:217:HIS:HB2	2.03	0.57
1:A:96:ARG:HA	1:A:182:TYR:O	2.06	0.56
1:A:155:ASP:CG	1:A:158:LYS:HB2	2.31	0.55
1:A:173:ASP:O	1:A:175:SER:HB3	2.05	0.55
1:A:119:LEU:C	1:A:119:LEU:HD13	2.32	0.54
1:A:83:PHE:HB2	1:A:194:LEU:O	2.07	0.54
1:A:155:ASP:OD2	1:A:162:LYS:HE2	2.07	0.54
1:A:82:ASP:OD2	1:A:85:LYS:HG3	2.08	0.54
1:A:113:LYS:HE3	2:A:351:HOH:O	2.08	0.54
1:A:103:ASP:OD2	1:A:131:LYS:HE3	2.08	0.54
1:A:151:TYR:HD2	1:A:200:TYR:CD1	2.26	0.54
1:A:133:ASP:OD1	1:A:133:ASP:N	2.40	0.54
1:A:171:ILE:HG22	1:A:172:GLU:N	2.22	0.53
1:A:178:LEU:CD2	1:A:179:ALA:H	2.21	0.53
1:A:173:ASP:O	1:A:175:SER:N	2.26	0.53
1:A:74:TYR:O	1:A:79:LYS:HE2	2.07	0.53
1:A:103:ASP:CG	1:A:131:LYS:HD2	2.34	0.53
1:A:151:TYR:HD2	1:A:200:TYR:HD1	1.57	0.53
1:A:173:ASP:O	1:A:173:ASP:OD1	2.26	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:ASP:CG	1:A:26:LYS:HG2	2.34	0.52
1:A:172:GLU:O	1:A:174:GLY:N	2.43	0.52
1:A:165:PHE:O	1:A:181:HIS:HB2	2.10	0.52
1:A:154:ALA:HA	1:A:161:ILE:HG22	1.92	0.52
1:A:110:ALA:HB2	1:A:123:ILE:HG12	1.91	0.51
1:A:44:LEU:HB3	1:A:46:PHE:CE1	2.45	0.51
1:A:100:PHE:CD2	1:A:136:ILE:HD11	2.46	0.51
1:A:122:ARG:NH2	2:A:346:HOH:O	2.44	0.51
1:A:165:PHE:O	1:A:181:HIS:N	2.32	0.50
1:A:103:ASP:HA	1:A:131:LYS:CE	2.41	0.50
1:A:18:LEU:HD13	1:A:19:ASP:N	2.26	0.50
1:A:78:MET:HE2	1:A:229:ILE:H	1.76	0.50
1:A:4:GLY:C	1:A:7:LEU:HD13	2.38	0.49
1:A:75:PRO:O	1:A:79:LYS:HG2	2.12	0.49
1:A:80:ARG:HG3	2:A:317:HOH:O	2.13	0.49
1:A:4:GLY:CA	1:A:7:LEU:HD13	2.42	0.48
1:A:171:ILE:HG22	1:A:173:ASP:N	2.27	0.48
1:A:61:VAL:O	1:A:66:CR2:HA11	2.13	0.48
1:A:135:ASN:CA	1:A:140:LYS:HG2	2.36	0.48
1:A:56:PRO:HD2	1:A:136:ILE:HG23	1.95	0.48
1:A:74:TYR:O	1:A:79:LYS:HD3	2.13	0.48
1:A:6:GLU:H	1:A:6:GLU:HG3	1.43	0.47
1:A:4:GLY:HA2	1:A:7:LEU:CD1	2.45	0.47
1:A:96:ARG:NE	1:A:183:GLN:OE1	2.48	0.47
1:A:229:ILE:O	1:A:230:THR:OG1	2.29	0.47
1:A:53:LEU:HD21	1:A:60:LEU:CD1	2.45	0.47
1:A:81:HIS:CD2	1:A:229:ILE:HG12	2.50	0.47
1:A:123:ILE:HG22	1:A:124:GLU:N	2.29	0.47
1:A:175:SER:OG	1:A:176:VAL:N	2.47	0.47
1:A:147:SER:OG	1:A:148:GLN:N	2.47	0.46
1:A:149:ASN:ND2	1:A:202:SER:HB2	2.30	0.46
1:A:149:ASN:HD22	1:A:202:SER:HB2	1.80	0.46
1:A:221:LEU:HD13	1:A:222:GLU:N	2.29	0.46
1:A:53:LEU:HD21	1:A:60:LEU:HD12	1.98	0.46
1:A:85:LYS:C	1:A:87:ALA:H	2.23	0.46
1:A:166:LYS:HA	1:A:179:ALA:O	2.15	0.46
1:A:201:LEU:HD13	1:A:203:TYR:HE1	1.80	0.45
1:A:148:GLN:CG	1:A:149:ASN:N	2.80	0.45
1:A:155:ASP:HB2	1:A:162:LYS:CG	2.40	0.45
1:A:8:PHE:CZ	1:A:88:MET:HG3	2.52	0.45
1:A:42:LEU:HD23	1:A:42:LEU:HA	1.79	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:PHE:CE2	1:A:136:ILE:HD13	2.51	0.45
1:A:151:TYR:CD2	1:A:200:TYR:HD1	2.31	0.44
1:A:153:MET:HE3	1:A:153:MET:HB3	1.58	0.44
1:A:160:GLY:C	1:A:161:ILE:HG23	2.43	0.44
1:A:18:LEU:HD13	1:A:19:ASP:O	2.18	0.43
1:A:181:HIS:N	1:A:181:HIS:ND1	2.66	0.43
1:A:51:GLY:C	1:A:52:LYS:HE3	2.43	0.43
1:A:150:VAL:HB	1:A:201:LEU:HD12	2.00	0.43
1:A:186:THR:OG1	1:A:187:PRO:HD2	2.18	0.43
1:A:78:MET:HE1	1:A:229:ILE:H	1.83	0.43
1:A:38:THR:O	1:A:73:ARG:HD3	2.18	0.43
1:A:78:MET:HE1	1:A:227:ALA:C	2.44	0.43
1:A:150:VAL:CG1	1:A:201:LEU:CD1	2.97	0.43
1:A:195:LEU:N	1:A:195:LEU:CD2	2.80	0.43
1:A:59:THR:HG22	1:A:100:PHE:CE2	2.54	0.43
1:A:110:ALA:HB2	1:A:123:ILE:CG1	2.48	0.43
1:A:73:ARG:O	1:A:73:ARG:HG2	2.18	0.42
1:A:130:PHE:HB3	1:A:137:LEU:HD12	2.01	0.42
1:A:178:LEU:CD2	1:A:179:ALA:N	2.80	0.42
1:A:95:GLU:O	1:A:183:GLN:HA	2.19	0.42
1:A:221:LEU:CD1	1:A:223:PHE:CE2	2.97	0.42
1:A:21:ASP:CB	1:A:26:LYS:HG2	2.49	0.42
1:A:176:VAL:CG1	1:A:177:GLN:N	2.79	0.42
1:A:18:LEU:HD13	1:A:18:LEU:C	2.44	0.42
1:A:55:VAL:HB	1:A:56:PRO:CD	2.48	0.42
1:A:194:LEU:C	1:A:195:LEU:HD22	2.44	0.42
1:A:178:LEU:HD23	1:A:178:LEU:HA	1.48	0.41
1:A:103:ASP:OD2	1:A:104:GLY:N	2.48	0.41
1:A:88:MET:HB3	1:A:89:PRO:HA	2.01	0.41
1:A:171:ILE:CG2	1:A:172:GLU:N	2.83	0.41
1:A:165:PHE:CD1	1:A:165:PHE:C	2.98	0.41
1:A:229:ILE:HD12	1:A:229:ILE:HA	1.60	0.41
1:A:123:ILE:CG2	1:A:124:GLU:N	2.83	0.41
1:A:140:LYS:N	1:A:140:LYS:HD2	2.35	0.41
1:A:18:LEU:C	1:A:18:LEU:CD1	2.93	0.41
1:A:42:LEU:HG	2:A:425:HOH:O	2.21	0.41
1:A:210:ASP:OD1	1:A:211:PRO:HD2	2.21	0.41
1:A:81:HIS:ND1	1:A:196:PRO:HA	2.36	0.40
1:A:4:GLY:CA	1:A:7:LEU:CD1	2.99	0.40
1:A:98:ILE:HG12	1:A:181:HIS:CD2	2.56	0.40
1:A:103:ASP:OD2	1:A:131:LYS:HD2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:LYS:CG	1:A:178:LEU:HD21	2.39	0.40
1:A:82:ASP:O	1:A:86:SER:HB3	2.21	0.40
1:A:78:MET:HE2	1:A:229:ILE:N	2.36	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	221/236 (94%)	204 (92%)	12 (5%)	5 (2%)	5 2

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	101	LYS
1	A	174	GLY
1	A	134	GLY
1	A	136	ILE
1	A	4	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	196/205 (96%)	148 (76%)	48 (24%)	1 0

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	5	GLU
1	A	6	GLU
1	A	7	LEU
1	A	9	THR
1	A	15	LEU
1	A	18	LEU
1	A	34	GLU
1	A	44	LEU
1	A	52	LYS
1	A	80	ARG
1	A	94	GLN
1	A	95	GLU
1	A	107	LYS
1	A	109	ARG
1	A	111	GLU
1	A	112	VAL
1	A	113	LYS
1	A	125	LEU
1	A	131	LYS
1	A	132	GLU
1	A	133	ASP
1	A	140	LYS
1	A	151	TYR
1	A	153	MET
1	A	156	LYS
1	A	157	GLN
1	A	158	LYS
1	A	161	ILE
1	A	166	LYS
1	A	167	ILE
1	A	168	ARG
1	A	170	ASN
1	A	173	ASP
1	A	178	LEU
1	A	180	ASP
1	A	186	THR
1	A	188	ILE
1	A	190	ASP
1	A	194	LEU
1	A	196	PRO
1	A	198	ASN

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Mol	Chain	Res	Type
1	A	201	LEU
1	A	202	SER
1	A	204	GLN
1	A	221	LEU
1	A	229	ILE
1	A	230	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	25	HIS
1	A	77	HIS
1	A	105	ASN
1	A	139	HIS
1	A	146	ASN
1	A	149	ASN
1	A	170	ASN
1	A	198	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	CR2	A	66	1	20,20,21	3.11	9 (45%)	25,27,29	2.30	9 (36%)

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	CR2	A	66	1	-	0/6/25/26	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	CR2	C1-N3	7.11	1.48	1.37
1	A	66	CR2	CG2-CB2	-6.02	1.35	1.46
1	A	66	CR2	CE1-CZ	5.15	1.48	1.39
1	A	66	CR2	OH-CZ	-4.77	1.26	1.37
1	A	66	CR2	CA1-C1	4.43	1.54	1.49
1	A	66	CR2	CA3-N3	-3.04	1.41	1.47
1	A	66	CR2	CE2-CZ	2.88	1.44	1.39
1	A	66	CR2	CD1-CG2	2.68	1.44	1.39
1	A	66	CR2	CE2-CD2	-2.00	1.35	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CR2	C2-N3-C1	-5.51	105.60	108.08
1	A	66	CR2	C3-CA3-N3	4.86	123.48	112.43
1	A	66	CR2	CA1-C1-N3	3.91	127.76	122.52
1	A	66	CR2	CG2-CB2-CA2	3.63	134.18	129.87
1	A	66	CR2	CB2-CA2-C2	-3.60	117.99	122.36
1	A	66	CR2	CE2-CZ-CE1	-2.46	115.84	119.77
1	A	66	CR2	C2-CA2-N2	2.38	110.66	108.95
1	A	66	CR2	CA2-C2-N3	2.31	105.44	103.50
1	A	66	CR2	CA1-C1-N2	-2.16	121.39	124.28

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
1	A	66	CR2	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	225/236 (95%)	0.18	2 (0%) 81 83	20, 32, 60, 87	11 (4%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	229	ILE	3.7
1	A	200	TYR	2.3

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	CR2	A	66	19/20	0.92	0.09	12,26,53,57	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.