



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 01:08 PM UTC

PDB ID : 8WGP / pdb_00008wgp
Title : Crystal structure of DsRed-Monomer
Authors : Nam, K.H.
Deposited on : 2023-09-22
Resolution : 2.90 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

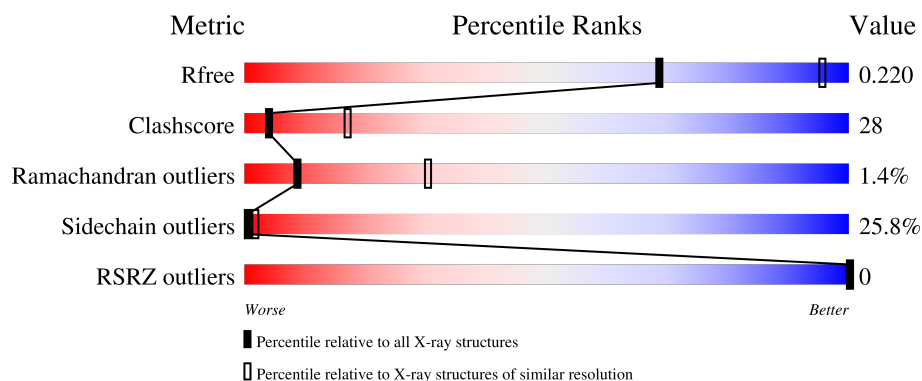
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

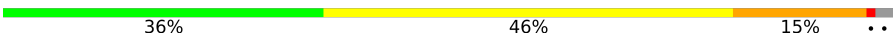
The reported resolution of this entry is 2.90 Å.

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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	223	 36% 46% 15% ..
1	B	223	 36% 49% 12% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3558 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 Red fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	0	2	0
			1755	1123	288	338	6			
1	B	218	Total	C	N	O	S	0	3	0
			1759	1126	288	339	6			

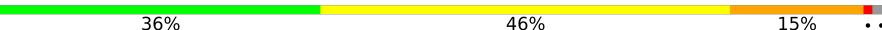
- Molecule 2 is water.

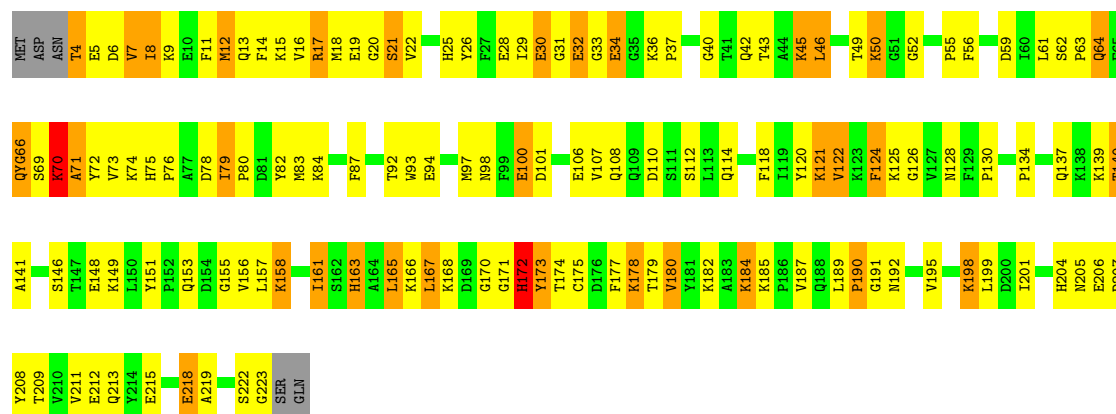
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	26	Total	O	0	0
			26	26		
2	B	18	Total	O	0	0
			18	18		

3 Residue-property plots

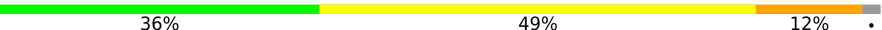
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

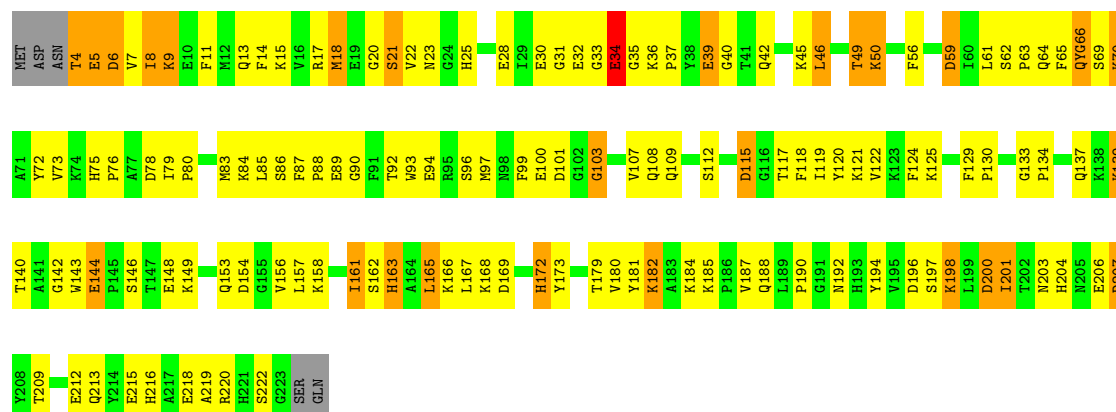
• Molecule 1: Red fluorescent protein

Chain A: 



• Molecule 1: Red fluorescent protein

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.57Å 107.00Å 50.03Å 90.00° 104.87° 90.00°	Depositor
Resolution (Å)	26.10 – 2.90 26.10 – 2.90	Depositor EDS
% Data completeness (in resolution range)	90.7 (26.10-2.90) 90.7 (26.10-2.90)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.36 (at 2.89Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.197 , 0.242 0.203 , 0.220	Depositor DCC
R_{free} test set	388 reflections (4.33%)	wwPDB-VP
Wilson B-factor (Å ²)	19.0	Xtriage
Anisotropy	0.509	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3558	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.20% 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: CRQ

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	1.10	0/1779	1.42	9/2402 (0.4%)
1	B	1.02	0/1786	1.38	9/2412 (0.4%)
All	All	1.06	0/3565	1.40	18/4814 (0.4%)

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

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	172[A]	HIS	CA-CB-CG	19.76	133.56	113.80
1	A	172[A]	HIS	CA-CB-CG	14.49	128.29	113.80
1	A	173	TYR	CB-CA-C	6.86	120.70	110.14
1	A	172[A]	HIS	CB-CA-C	6.15	120.59	109.83
1	A	36	LYS	CB-CA-C	6.02	115.86	110.44
1	B	59	ASP	CB-CA-C	5.95	120.66	110.79
1	A	71	ALA	N-CA-C	-5.90	105.84	113.16
1	A	55	PRO	N-CA-CB	-5.48	99.14	102.81
1	B	103	GLY	CA-C-O	-5.47	116.60	121.41
1	A	126	GLY	CA-C-O	-5.46	116.33	120.76
1	B	142	GLY	CA-C-O	-5.42	117.52	122.24
1	B	42	GLN	CA-C-O	-5.34	115.36	121.40
1	B	117	THR	CA-CB-OG1	-5.32	101.62	109.60
1	B	73	VAL	CA-C-O	-5.18	115.04	120.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	179	THR	CA-CB-OG1	-5.13	101.90	109.60
1	A	70	LYS	N-CA-C	-5.05	107.16	113.38
1	B	144[A]	GLU	CB-CG-CD	5.05	121.19	112.60
1	A	191	GLY	CA-C-O	-5.01	115.84	121.31

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	52	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	1755	0	1660	104	4
1	B	1759	0	1666	89	6
2	A	26	0	0	4	0
2	B	18	0	0	4	0
All	All	3558	0	3326	190	6

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 28.

All (190) 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:30:GLU:HG3	1:A:49:THR:HG21	1.52	0.92
1:B:144[A]:GLU:HG3	1:B:166:LYS:HG2	1.53	0.90
1:A:153:GLN:HB3	1:A:158:LYS:HE3	1.58	0.86
1:A:31:GLY:HA3	1:A:46:LEU:HD12	1.58	0.85
1:A:149:LYS:HG3	2:A:317:HOH:O	1.76	0.84
1:A:100:GLU:OE2	1:A:174:THR:N	2.10	0.84
1:A:167:LEU:HB2	1:A:171:GLY:O	1.81	0.81
1:B:134:PRO:HD3	2:B:310:HOH:O	1.78	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:ILE:HD12	1:B:84:LYS:HG2	1.65	0.78
1:A:134:PRO:HB3	1:A:167:LEU:HD21	1.66	0.77
1:B:76:PRO:HG3	2:B:309:HOH:O	1.84	0.77
1:A:4:THR:HG21	1:A:84:LYS:O	1.85	0.76
1:A:101:ASP:HB2	1:A:130:PRO:HG2	1.69	0.74
1:B:207:ASP:OD2	1:B:209:THR:OG1	2.04	0.74
1:A:70:LYS:HB2	1:A:83:MET:HG2	1.69	0.74
1:A:134:PRO:CB	1:A:167:LEU:HD21	2.19	0.72
1:B:133:GLY:HA2	2:B:310:HOH:O	1.87	0.72
1:A:161:ILE:HB	1:A:177:PHE:HB2	1.71	0.72
1:A:42:GLN:HE22	1:A:69:SER:HB3	1.55	0.72
1:B:56:PHE:HE1	1:B:61:LEU:HD11	1.56	0.71
1:A:17:ARG:NH1	1:A:19:GLU:OE2	2.23	0.70
1:B:13:GLN:HG2	1:B:34:GLU:HB3	1.74	0.70
1:A:146:SER:HB3	1:A:163:HIS:CE1	2.26	0.69
1:B:4:THR:HG21	1:B:84:LYS:HB3	1.76	0.68
1:B:22:VAL:HG22	1:B:124:PHE:HE1	1.59	0.67
1:A:42:GLN:NE2	1:A:69:SER:HB3	2.10	0.67
1:A:75:HIS:HA	1:A:219:ALA:HB3	1.75	0.66
1:A:7:VAL:HG12	1:A:8:ILE:HG12	1.78	0.66
1:A:18:MET:HG3	1:A:122:VAL:HB	1.77	0.66
1:B:134:PRO:HA	1:B:139:LYS:HG3	1.80	0.64
1:A:166:LYS:NZ	1:B:200:ASP:OD2	2.28	0.64
1:A:139:LYS:O	1:A:168:LYS:HG3	1.97	0.64
1:B:207:ASP:HB2	1:B:209:THR:HG23	1.80	0.63
1:B:94:GLU:HG2	1:B:108:GLN:HG2	1.79	0.63
1:A:30:GLU:CG	1:A:49:THR:HG21	2.27	0.62
1:B:25:HIS:HB2	2:B:311:HOH:O	1.98	0.62
1:B:137:GLN:HB3	1:B:139:LYS:HD3	1.80	0.62
1:B:101:ASP:OD2	1:B:173:TYR:OH	2.18	0.62
1:A:170:GLY:O	1:B:203:ASN:OD1	2.17	0.62
1:A:76:PRO:HD2	1:A:79:ILE:HB	1.82	0.60
1:B:23:ASN:HD21	1:B:129:PHE:H	1.49	0.60
1:A:112:SER:OG	1:A:114:GLN:OE1	2.17	0.60
1:A:74:LYS:HD2	1:A:218:GLU:CD	2.26	0.60
1:A:16:VAL:HG23	1:A:120:TYR:HB2	1.84	0.60
1:A:22:VAL:HG21	1:A:56:PHE:CZ	2.36	0.60
1:B:75:HIS:HA	1:B:219:ALA:HB3	1.84	0.60
1:A:28:GLU:HB2	1:A:50:LYS:HB3	1.84	0.59
1:A:82:TYR:HB2	1:A:190:PRO:HD3	1.84	0.59
1:B:109:GLN:HG3	1:B:122:VAL:HG22	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:LYS:HG2	1:A:212:GLU:HG2	1.84	0.59
1:B:201:ILE:N	1:B:201:ILE:HD13	2.16	0.59
1:A:42:GLN:HE22	1:A:69:SER:CB	2.15	0.59
1:B:156:VAL:HG12	1:B:182:LYS:HG3	1.85	0.59
1:A:205:ASN:ND2	1:A:207:ASP:OD1	2.35	0.59
1:B:83:MET:HE3	1:B:181:TYR:CD1	2.38	0.58
1:B:101:ASP:CG	1:B:173:TYR:OH	2.46	0.58
1:A:97:MET:HE2	1:A:177:PHE:CZ	2.39	0.58
1:A:40:GLY:HA2	1:A:72:TYR:O	2.03	0.57
1:B:153:GLN:O	1:B:156:VAL:HG22	2.03	0.57
1:A:7:VAL:HG11	1:A:87:PHE:HB3	1.85	0.57
1:A:97:MET:HE2	1:A:177:PHE:CE2	2.40	0.57
1:B:90:GLY:C	1:B:184:LYS:HG3	2.29	0.57
1:B:22:VAL:HG22	1:B:124:PHE:CE1	2.39	0.57
1:B:146:SER:HB3	1:B:163:HIS:CE1	2.40	0.57
1:B:21:SER:HA	1:B:25:HIS:O	2.05	0.56
1:B:18:MET:HE3	1:B:65:PHE:CZ	2.40	0.56
1:A:74:LYS:HD2	1:A:218:GLU:OE1	2.05	0.56
1:B:139:LYS:HA	1:B:168:LYS:HD2	1.87	0.56
1:B:8:ILE:HD13	1:B:118:PHE:HZ	1.72	0.55
1:B:36:LYS:HB3	1:B:39:GLU:HG3	1.87	0.55
1:A:195:VAL:HG22	1:A:219:ALA:HA	1.88	0.55
1:A:128:ASN:O	1:A:130:PRO:HD3	2.07	0.54
1:B:31:GLY:HA3	1:B:46:LEU:HD12	1.89	0.54
1:A:13:GLN:HA	1:A:34:GLU:HA	1.90	0.54
1:A:70:LYS:HA	1:A:73:VAL:HG23	1.89	0.54
1:A:100:GLU:OE2	1:A:173:TYR:HA	2.08	0.54
1:A:155:GLY:C	1:A:189:LEU:HD11	2.33	0.53
1:A:62:SER:OG	1:A:63:PRO:HD3	2.07	0.53
1:B:35:GLY:O	1:B:37:PRO:HD3	2.08	0.53
1:A:11:PHE:HE1	1:A:34:GLU:HB2	1.74	0.53
1:A:66:CRQ:HE11	1:A:215:GLU:HB2	1.73	0.53
1:A:161:ILE:HG22	1:A:163:HIS:HB2	1.91	0.53
1:A:93:TRP:HA	1:A:180:VAL:O	2.08	0.52
1:B:201:ILE:HD12	1:B:213:GLN:HG2	1.91	0.52
1:A:137:GLN:HB2	1:A:139:LYS:HD3	1.91	0.52
1:B:207:ASP:OD1	1:B:207:ASP:N	2.42	0.52
1:A:49:THR:O	1:A:50:LYS:HB2	2.09	0.52
1:B:93:TRP:HA	1:B:180:VAL:O	2.10	0.52
1:A:45:LYS:CG	1:A:212:GLU:HG2	2.40	0.51
1:A:37:PRO:HB3	1:A:71:ALA:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:THR:HG21	1:A:165:LEU:HD13	1.93	0.51
1:B:31:GLY:HA3	1:B:46:LEU:CD1	2.41	0.51
1:A:161:ILE:CG2	1:A:163:HIS:HB2	2.40	0.51
1:A:61:LEU:O	1:A:64:GLN:HG2	2.11	0.51
1:B:157:LEU:HD13	1:B:190:PRO:HG2	1.91	0.51
1:B:45:LYS:HE3	1:B:212:GLU:OE1	2.10	0.51
1:A:97:MET:HB3	1:A:175:CYS:SG	2.51	0.51
1:B:148:GLU:HB2	1:B:161:ILE:HG12	1.92	0.50
1:B:66:CRQ:HB1	1:B:70:LYS:HE3	1.94	0.50
1:A:15:LYS:HG2	1:A:32:GLU:HG3	1.94	0.50
1:B:66:CRQ:OH	1:B:146:SER:OG	2.29	0.50
1:A:14:PHE:CZ	1:A:33:GLY:HA3	2.47	0.50
1:B:76:PRO:HD2	1:B:79:ILE:HB	1.94	0.50
1:A:45:LYS:HE3	1:A:212:GLU:OE1	2.12	0.49
1:B:11:PHE:CZ	1:B:34:GLU:HB2	2.47	0.49
1:B:15:LYS:HG2	1:B:32:GLU:HG3	1.95	0.49
1:B:87:PHE:HB3	1:B:88:PRO:HA	1.94	0.49
1:A:134:PRO:HB2	1:A:167:LEU:HD21	1.94	0.49
1:B:20:GLY:HA2	1:B:124:PHE:O	2.12	0.49
1:B:59:ASP:O	1:B:63:PRO:HD3	2.12	0.49
1:B:62:SER:OG	1:B:63:PRO:HD3	2.11	0.49
1:B:70:LYS:CB	1:B:83:MET:HE2	2.43	0.49
1:A:14:PHE:HB3	1:A:118:PHE:HB2	1.95	0.49
1:A:37:PRO:HA	1:A:72:TYR:HA	1.95	0.49
1:A:14:PHE:HA	1:A:118:PHE:O	2.13	0.48
1:B:156:VAL:HG12	1:B:182:LYS:CG	2.43	0.48
1:B:86:SER:O	1:B:90:GLY:N	2.33	0.48
1:A:201:ILE:HD12	1:A:213:GLN:HG2	1.96	0.47
1:A:64:GLN:HE22	1:A:97:MET:CE	2.27	0.47
1:A:21:SER:HA	1:A:25:HIS:O	2.15	0.47
1:B:49:THR:O	1:B:50:LYS:HB2	2.15	0.47
1:B:59:ASP:HB3	1:B:165:LEU:HD21	1.97	0.46
1:A:20:GLY:HA2	1:A:124:PHE:O	2.15	0.46
1:B:101:ASP:HB2	1:B:130:PRO:HG2	1.96	0.46
1:A:66:CRQ:NE1	1:A:199:LEU:HD11	2.30	0.46
1:B:137:GLN:CB	1:B:139:LYS:HD3	2.46	0.46
1:A:110:ASP:OD1	1:A:184:LYS:HE2	2.14	0.46
1:B:143:TRP:CD2	1:B:165:LEU:HG	2.51	0.46
1:B:169:ASP:C	1:B:169:ASP:OD1	2.57	0.46
1:B:85:LEU:HB2	1:B:187:VAL:HG21	1.98	0.46
1:A:50:LYS:HA	2:A:320:HOH:O	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LYS:HE2	1:A:151:TYR:CE1	2.50	0.45
1:B:13:GLN:HA	1:B:34:GLU:HA	1.98	0.45
1:B:203:ASN:O	1:B:204:HIS:HB3	2.17	0.45
1:B:28:GLU:HB2	1:B:50:LYS:HB3	2.00	0.44
1:B:167:LEU:HD23	1:B:167:LEU:HA	1.86	0.44
1:A:205:ASN:ND2	1:A:209:THR:OG1	2.46	0.44
1:A:22:VAL:HG21	1:A:56:PHE:CE1	2.53	0.44
1:A:76:PRO:HD3	1:A:219:ALA:HB3	2.00	0.44
1:B:40:GLY:O	1:B:216:HIS:HA	2.17	0.44
1:B:23:ASN:ND2	1:B:129:PHE:HB2	2.32	0.44
1:A:204:HIS:CE1	1:A:208:TYR:CZ	3.06	0.44
1:B:101:ASP:CG	1:B:173:TYR:HH	2.20	0.44
1:A:157:LEU:HD13	1:A:190:PRO:HG2	2.00	0.43
1:B:18:MET:HE1	1:B:61:LEU:HD22	1.99	0.43
1:A:110:ASP:HB3	1:A:121:LYS:HE3	1.99	0.43
1:B:203:ASN:HB3	1:B:212:GLU:HB2	2.00	0.43
1:A:141:ALA:HB2	1:A:168:LYS:HG2	2.00	0.43
1:B:17:ARG:NH2	1:B:28:GLU:OE1	2.49	0.43
1:A:12:MET:HE3	1:A:12:MET:HB3	1.68	0.43
1:A:100:GLU:OE2	1:A:174:THR:HG23	2.19	0.43
1:B:9:LYS:H	1:B:9:LYS:HG2	1.65	0.43
1:B:45:LYS:O	1:B:46:LEU:HD13	2.19	0.43
1:B:37:PRO:HG3	1:B:72:TYR:CE2	2.54	0.42
1:B:109:GLN:HE21	1:B:120:TYR:HB3	1.83	0.42
1:A:14:PHE:CB	1:A:118:PHE:HB2	2.50	0.42
1:B:18:MET:HG3	1:B:122:VAL:HB	2.00	0.42
1:A:12:MET:HE2	1:A:118:PHE:CE2	2.55	0.42
1:A:180:VAL:O	1:A:180:VAL:HG12	2.18	0.42
1:A:66:CRQ:HD2	1:A:215:GLU:OE1	2.19	0.42
1:A:198:LYS:HE2	2:A:314:HOH:O	2.19	0.42
1:A:26:TYR:CE2	1:A:28:GLU:HG3	2.55	0.42
1:A:101:ASP:OD1	1:A:173:TYR:OH	2.29	0.42
1:B:23:ASN:HD21	1:B:129:PHE:HB2	1.84	0.42
1:A:18:MET:HA	1:A:122:VAL:O	2.20	0.42
1:A:204:HIS:HA	1:A:211:VAL:HA	2.02	0.42
1:A:98:ASN:HA	2:A:306:HOH:O	2.19	0.41
1:A:158:LYS:HB3	1:A:178:LYS:HE2	2.01	0.41
1:B:28:GLU:O	1:B:49:THR:N	2.45	0.41
1:A:156:VAL:N	1:A:189:LEU:HD11	2.35	0.41
1:A:172[A]:HIS:CE1	1:A:174:THR:HG22	2.55	0.41
1:A:223:GLY:O	1:B:36:LYS:NZ	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:MET:HE3	1:B:65:PHE:CE1	2.55	0.41
1:B:197:SER:HA	1:B:216:HIS:O	2.20	0.41
1:A:59:ASP:HB3	1:A:165:LEU:HD21	2.02	0.41
1:B:17:ARG:HH22	1:B:28:GLU:CD	2.28	0.41
1:B:134:PRO:HB3	1:B:167:LEU:HD21	2.02	0.41
1:B:198:LYS:O	1:B:215:GLU:HA	2.21	0.41
1:A:79:ILE:HA	1:A:80:PRO:HD3	1.85	0.41
1:B:79:ILE:HA	1:B:80:PRO:HD3	1.87	0.41
1:A:18:MET:CE	1:A:64:GLN:HB2	2.51	0.41
1:A:124:PHE:CD1	1:A:124:PHE:C	2.98	0.41
1:B:4:THR:O	1:B:6:ASP:N	2.53	0.41
1:B:14:PHE:CZ	1:B:33:GLY:HA3	2.55	0.41
1:A:16:VAL:HA	1:A:120:TYR:O	2.20	0.40
1:A:66:CRQ:NE1	1:A:215:GLU:HB2	2.36	0.40
1:B:103:GLY:HA3	1:B:129:PHE:CE1	2.56	0.40
1:A:148:GLU:HG3	1:A:161:ILE:HG13	2.04	0.40
1:A:94:GLU:HG2	1:A:108:GLN:HG2	2.04	0.40

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:LYS:NZ	1:B:154:ASP:OD2[2_454]	1.05	1.15
1:A:34:GLU:OE1	1:B:34:GLU:OE1[1_455]	1.44	0.76
1:A:45:LYS:NZ	1:B:45:LYS:NZ[1_455]	1.98	0.22
1:B:115:ASP:OD1	1:B:194:TYR:CE2[1_655]	2.11	0.09
1:A:50:LYS:NZ	1:B:154:ASP:CG[2_454]	2.14	0.06
1:B:115:ASP:OD2	1:B:220:ARG:NH2[1_655]	2.19	0.01

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	213/223 (96%)	195 (92%)	15 (7%)	3 (1%)	9	30
1	B	214/223 (96%)	193 (90%)	18 (8%)	3 (1%)	9	30
All	All	427/446 (96%)	388 (91%)	33 (8%)	6 (1%)	9	30

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	50	LYS
1	A	190	PRO
1	B	50	LYS
1	B	5	GLU
1	A	192	ASN
1	B	34	GLU

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	186/191 (97%)	140 (75%)	46 (25%)	1	2
1	B	187/191 (98%)	137 (73%)	50 (27%)	0	1
All	All	373/382 (98%)	277 (74%)	96 (26%)	0	2

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	5	GLU
1	A	6	ASP
1	A	7	VAL
1	A	8	ILE
1	A	9	LYS
1	A	12	MET
1	A	17	ARG
1	A	21	SER
1	A	29	ILE

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Mol	Chain	Res	Type
1	A	30	GLU
1	A	32	GLU
1	A	34	GLU
1	A	43	THR
1	A	45	LYS
1	A	46	LEU
1	A	64	GLN
1	A	70	LYS
1	A	78	ASP
1	A	79	ILE
1	A	92	THR
1	A	100	GLU
1	A	106	GLU
1	A	107	VAL
1	A	121	LYS
1	A	122	VAL
1	A	124	PHE
1	A	125	LYS
1	A	140	THR
1	A	158	LYS
1	A	161	ILE
1	A	163	HIS
1	A	165	LEU
1	A	167	LEU
1	A	172[A]	HIS
1	A	178	LYS
1	A	179	THR
1	A	180	VAL
1	A	182	LYS
1	A	184	LYS
1	A	185	LYS
1	A	187	VAL
1	A	198	LYS
1	A	206	GLU
1	A	218	GLU
1	A	222	SER
1	B	4	THR
1	B	5	GLU
1	B	6	ASP
1	B	7	VAL
1	B	8	ILE
1	B	9	LYS

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Mol	Chain	Res	Type
1	B	18	MET
1	B	21	SER
1	B	30	GLU
1	B	34	GLU
1	B	39	GLU
1	B	46	LEU
1	B	49	THR
1	B	64	GLN
1	B	69	SER
1	B	70	LYS
1	B	78	ASP
1	B	89	GLU
1	B	92	THR
1	B	96	SER
1	B	97	MET
1	B	99	PHE
1	B	100	GLU
1	B	107	VAL
1	B	112	SER
1	B	115	ASP
1	B	119	ILE
1	B	121	LYS
1	B	125	LYS
1	B	139	LYS
1	B	140	THR
1	B	149	LYS
1	B	158	LYS
1	B	161	ILE
1	B	162	SER
1	B	163	HIS
1	B	165	LEU
1	B	172[A]	HIS
1	B	182	LYS
1	B	185	LYS
1	B	188	GLN
1	B	192	ASN
1	B	196	ASP
1	B	198	LYS
1	B	200	ASP
1	B	201	ILE
1	B	206	GLU
1	B	207	ASP

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Mol	Chain	Res	Type
1	B	218	GLU
1	B	222	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	64	GLN
1	A	75	HIS
1	A	172[A]	HIS
1	B	23	ASN
1	B	192	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 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	CRQ	B	66	1	25,25,26	0.92	1 (4%)	28,34,36	1.08	2 (7%)
1	CRQ	A	66	1	25,25,26	0.94	1 (4%)	28,34,36	1.35	4 (14%)

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	CRQ	B	66	1	-	2/10/32/33	0/2/2/2
1	CRQ	A	66	1	-	5/10/32/33	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	66	CRQ	C1-N2	2.97	1.39	1.33
1	B	66	CRQ	C1-N2	2.55	1.38	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	CRQ	O3-C3-CA3	-4.13	106.87	125.47
1	B	66	CRQ	O3-C3-CA3	-3.90	107.93	125.47
1	A	66	CRQ	O2-C2-CA2	-3.40	128.85	131.02
1	B	66	CRQ	N3-C1-N2	-2.29	109.90	112.62
1	A	66	CRQ	C3-CA3-N3	2.27	117.59	112.43
1	A	66	CRQ	CA2-C2-N3	2.07	105.24	103.50

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	66	CRQ	C2-CA2-CB2-CG2
1	B	66	CRQ	C1-CA1-CB1-CG1
1	B	66	CRQ	CA1-CB1-CG1-CD3
1	A	66	CRQ	CA1-CB1-CG1-CD3
1	A	66	CRQ	N2-CA2-CB2-CG2
1	A	66	CRQ	OE1-CD3-CG1-CB1
1	A	66	CRQ	NE1-CD3-CG1-CB1

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	66	CRQ	2	0
1	A	66	CRQ	4	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

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	217/223 (97%)	0.01	0 100 100	3, 12, 23, 60	2 (0%)
1	B	217/223 (97%)	0.04	0 100 100	3, 12, 25, 46	3 (1%)
All	All	434/446 (97%)	0.03	0 100 100	3, 12, 25, 60	5 (1%)

There are no RSRZ outliers to report.

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	CRQ	A	66	24/25	0.92	0.13	8,24,52,70	0
1	CRQ	B	66	24/25	0.93	0.09	8,18,33,33	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.