



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:11 AM UTC

PDB ID : 5WHF / pdb_00005whf
Title : Crystal structure of vimentin coil 1B packed in a high-order filamentous form
Authors : Obiero, J.M.; Pang, A.H.; Tsodikov, O.V.
Deposited on : 2017-07-16
Resolution : 2.25 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

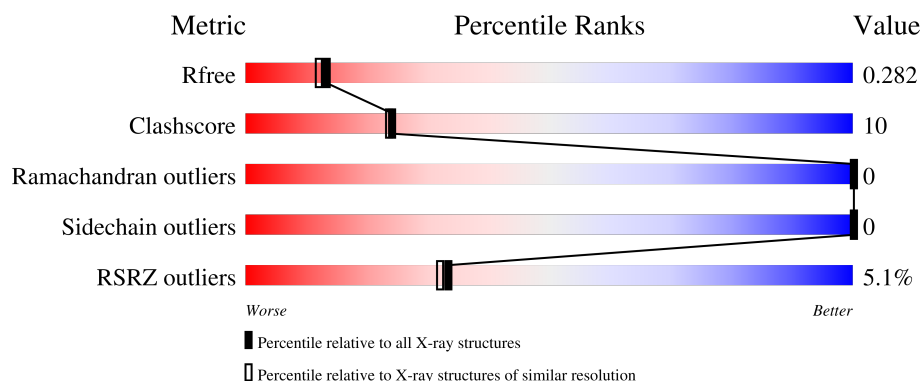
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

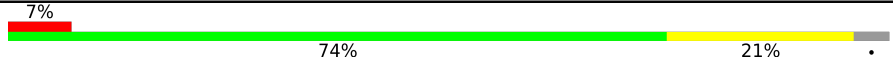
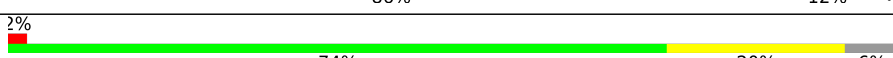
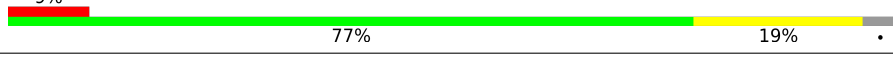
The reported resolution of this entry is 2.25 Å.

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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	90	
1	B	90	
1	C	90	
1	D	90	
1	E	90	

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Mol	Chain	Length	Quality of chain
1	F	90	2% 82% 11% 7%
1	G	90	6% 84% 13% 7%
1	H	90	2% 80% 13% 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5475 atoms, of which 8 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 Vimentin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	86	Total	C	N	O	S	0	0	0
			662	405	125	130	2			
1	B	82	Total	C	N	O	S	0	0	0
			647	395	122	128	2			
1	C	88	Total	C	N	O	S	0	0	0
			689	420	131	134	4			
1	D	85	Total	C	N	O	S	0	0	0
			682	416	128	135	3			
1	E	86	Total	C	N	O	S	0	0	0
			642	395	119	127	1			
1	F	84	Total	C	N	O	S	0	0	0
			641	390	121	127	3			
1	G	88	Total	C	N	O	S	0	0	0
			696	426	134	133	3			
1	H	84	Total	C	N	O	S	0	0	0
			675	410	129	133	3			

There are 32 discrepancies between the modelled and reference sequences:

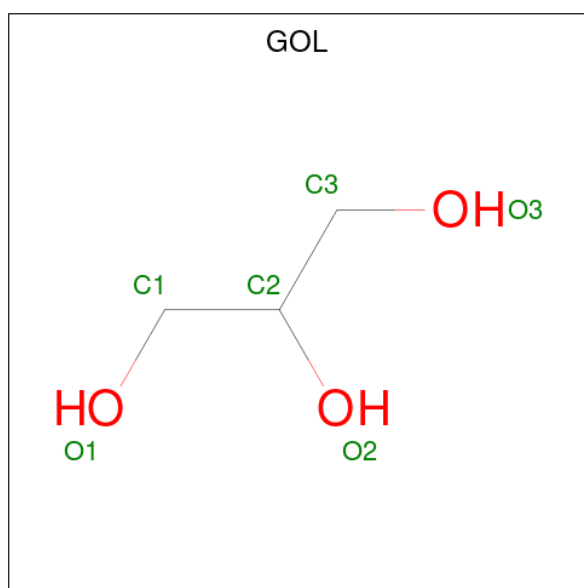
Chain	Residue	Modelled	Actual	Comment	Reference
A	149	GLY	-	expression tag	UNP P08670
A	150	PRO	-	expression tag	UNP P08670
A	151	HIS	-	expression tag	UNP P08670
A	152	MET	-	expression tag	UNP P08670
B	149	GLY	-	expression tag	UNP P08670
B	150	PRO	-	expression tag	UNP P08670
B	151	HIS	-	expression tag	UNP P08670
B	152	MET	-	expression tag	UNP P08670
C	149	GLY	-	expression tag	UNP P08670
C	150	PRO	-	expression tag	UNP P08670
C	151	HIS	-	expression tag	UNP P08670
C	152	MET	-	expression tag	UNP P08670
D	149	GLY	-	expression tag	UNP P08670

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Chain	Residue	Modelled	Actual	Comment	Reference
D	150	PRO	-	expression tag	UNP P08670
D	151	HIS	-	expression tag	UNP P08670
D	152	MET	-	expression tag	UNP P08670
E	149	GLY	-	expression tag	UNP P08670
E	150	PRO	-	expression tag	UNP P08670
E	151	HIS	-	expression tag	UNP P08670
E	152	MET	-	expression tag	UNP P08670
F	149	GLY	-	expression tag	UNP P08670
F	150	PRO	-	expression tag	UNP P08670
F	151	HIS	-	expression tag	UNP P08670
F	152	MET	-	expression tag	UNP P08670
G	149	GLY	-	expression tag	UNP P08670
G	150	PRO	-	expression tag	UNP P08670
G	151	HIS	-	expression tag	UNP P08670
G	152	MET	-	expression tag	UNP P08670
H	149	GLY	-	expression tag	UNP P08670
H	150	PRO	-	expression tag	UNP P08670
H	151	HIS	-	expression tag	UNP P08670
H	152	MET	-	expression tag	UNP P08670

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	G	1	Total	C	H	O	0	0
			14	3	8	3		

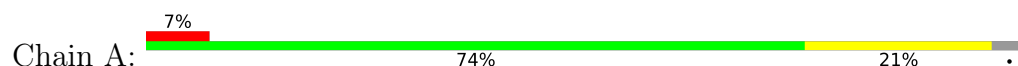
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total 15	O 15	0	0
3	B	7	Total 7	O 7	0	0
3	C	14	Total 14	O 14	0	0
3	D	20	Total 20	O 20	0	0
3	E	14	Total 14	O 14	0	0
3	F	19	Total 19	O 19	0	0
3	G	17	Total 17	O 17	0	0
3	H	21	Total 21	O 21	0	0

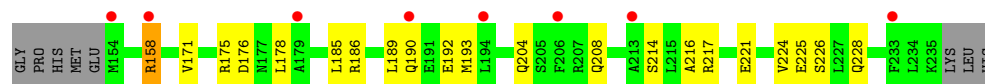
3 Residue-property plots [i](#)

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.

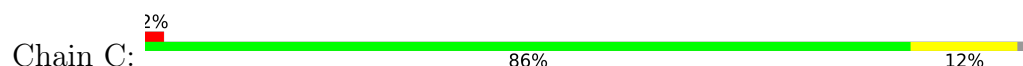
- Molecule 1: Vimentin



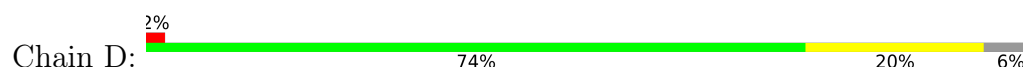
- Molecule 1: Vimentin



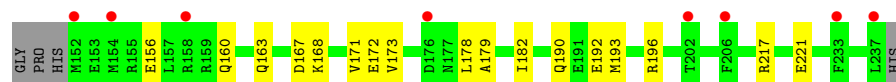
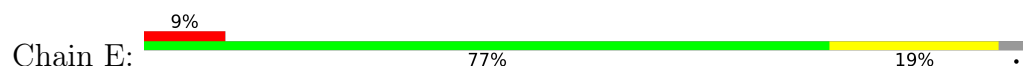
- Molecule 1: Vimentin



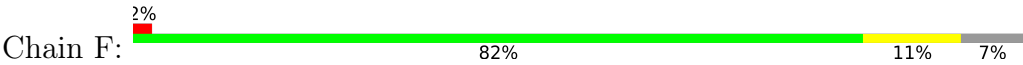
- Molecule 1: Vimentin



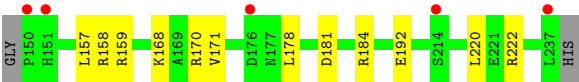
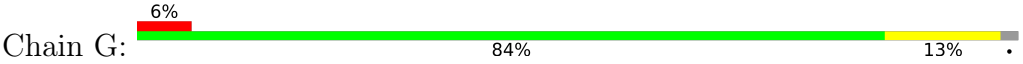
- Molecule 1: Vimentin



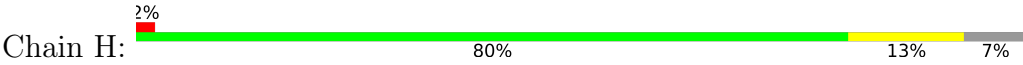
- Molecule 1: Vimentin



• Molecule 1: Vimentin



• Molecule 1: Vimentin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.99Å 77.27Å 123.39Å 90.00° 95.03° 90.00°	Depositor
Resolution (Å)	40.97 – 2.25 40.97 – 2.25	Depositor EDS
% Data completeness (in resolution range)	89.3 (40.97-2.25) 89.3 (40.97-2.25)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.95 (at 2.24Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.235 , 0.282 0.235 , 0.282	Depositor DCC
R_{free} test set	1601 reflections (3.60%)	wwPDB-VP
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5475	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.65 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.0745e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GOL

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	0.43	0/663	0.59	0/892
1	B	0.44	0/648	0.60	0/870
1	C	0.33	0/691	0.52	0/926
1	D	0.31	0/683	0.47	0/915
1	E	0.36	0/643	0.51	0/867
1	F	0.29	0/642	0.44	0/864
1	G	0.32	0/698	0.54	0/933
1	H	0.31	0/675	0.53	0/903
All	All	0.35	0/5343	0.53	0/7170

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
1	D	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	158	ARG	Sidechain
1	D	222	ARG	Sidechain

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	662	0	632	21	0
1	B	647	0	624	24	0
1	C	689	0	671	16	0
1	D	682	0	665	27	0
1	E	642	0	596	19	0
1	F	641	0	600	18	0
1	G	696	0	686	16	0
1	H	675	0	660	15	0
2	G	6	8	8	0	0
3	A	15	0	0	0	0
3	B	7	0	0	1	0
3	C	14	0	0	1	0
3	D	20	0	0	0	0
3	E	14	0	0	1	0
3	F	19	0	0	0	0
3	G	17	0	0	2	0
3	H	21	0	0	1	0
All	All	5467	8	5142	105	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 10.

All (105) 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:171:VAL:HG12	1:B:171:VAL:HG12	1.50	0.91
1:E:171:VAL:HG12	1:F:171:VAL:HG12	1.57	0.86
1:D:189:LEU:HD11	1:D:193:MET:HE3	1.57	0.84
1:B:189:LEU:HD11	1:B:193:MET:HE3	1.62	0.81
1:A:190:GLN:O	1:A:194:LEU:HD13	1.81	0.79
1:E:190:GLN:HA	1:E:193:MET:HE2	1.66	0.77
1:F:189:LEU:HD11	1:F:193:MET:HE3	1.70	0.73
1:E:173:VAL:HG23	3:E:301:HOH:O	1.89	0.73
1:E:179:ALA:O	1:E:182:ILE:HG22	1.89	0.72
1:C:178:LEU:HD22	1:D:182:ILE:HD12	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:LEU:CD1	1:D:217:ARG:HG3	2.19	0.72
1:F:217:ARG:NH1	1:F:217:ARG:HB2	2.05	0.71
1:G:171:VAL:HG12	1:H:171:VAL:HG12	1.74	0.68
1:C:171:VAL:HG12	1:D:171:VAL:HG12	1.76	0.67
1:E:168:LYS:NZ	1:E:172:GLU:OE2	2.24	0.67
1:C:175:ARG:HD2	1:D:174:GLU:CD	2.20	0.66
1:C:174:GLU:OE1	1:D:175:ARG:NH1	2.29	0.66
1:C:178:LEU:CD2	1:D:182:ILE:HD12	2.27	0.65
1:B:189:LEU:CD1	1:B:193:MET:HE3	2.27	0.65
3:C:301:HOH:O	1:D:193:MET:HE2	1.97	0.64
1:G:178:LEU:HD22	1:H:182:ILE:HD12	1.80	0.64
1:A:227:LEU:HD11	1:B:228:GLN:NE2	2.13	0.64
1:A:233:PHE:CE1	1:D:154:MET:HE1	2.33	0.63
1:C:220:LEU:HD11	1:D:217:ARG:HG3	1.80	0.62
1:F:162:ASP:OD1	1:G:222:ARG:NE	2.29	0.61
1:G:158:ARG:HG3	1:H:157:LEU:HD11	1.83	0.60
1:C:182:ILE:HG12	1:D:182:ILE:HG12	1.82	0.60
1:E:171:VAL:CG1	1:F:171:VAL:HG12	2.30	0.60
1:B:216:ALA:HA	3:B:301:HOH:O	2.03	0.59
1:E:193:MET:HG2	1:F:192:GLU:OE1	2.02	0.59
1:G:159:ARG:NH2	3:G:401:HOH:O	2.34	0.58
1:E:171:VAL:HG12	1:F:171:VAL:CG1	2.33	0.58
1:D:217:ARG:CZ	1:D:217:ARG:HB2	2.34	0.57
1:B:224:VAL:HG13	1:B:228:GLN:HE22	1.68	0.57
1:A:219:ASP:OD1	1:D:166:ASN:ND2	2.37	0.57
1:A:233:PHE:CZ	1:D:154:MET:HE1	2.40	0.57
1:D:182:ILE:O	1:D:186:ARG:HG3	2.05	0.57
1:B:175:ARG:HH21	1:B:176:ASP:CG	2.13	0.56
1:G:181:ASP:OD1	1:G:184:ARG:NH2	2.38	0.56
1:G:192:GLU:OE2	1:H:196:ARG:NE	2.33	0.56
1:F:189:LEU:CD1	1:F:193:MET:HE3	2.35	0.56
1:C:182:ILE:CG1	1:D:182:ILE:HG12	2.37	0.55
1:E:192:GLU:HG2	1:F:193:MET:HG3	1.90	0.53
1:G:170:ARG:HD2	3:G:402:HOH:O	2.07	0.53
1:C:154:MET:SD	1:D:154:MET:HG3	2.49	0.52
1:A:157:LEU:HD11	1:B:158:ARG:HH21	1.73	0.52
1:F:217:ARG:HB2	1:F:217:ARG:CZ	2.40	0.52
1:A:171:VAL:CG1	1:B:171:VAL:HG12	2.31	0.51
1:A:175:ARG:HG2	1:B:178:LEU:CD1	2.41	0.51
1:C:182:ILE:HD12	1:D:178:LEU:HD22	1.92	0.51
1:D:189:LEU:CD1	1:D:193:MET:HE3	2.36	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:193:MET:HG2	1:F:192:GLU:CD	2.36	0.50
1:A:161:VAL:O	1:A:165:THR:HG23	2.11	0.50
1:A:236:LYS:O	1:A:236:LYS:HG3	2.11	0.50
1:A:181:ASP:OD1	1:A:184:ARG:NH2	2.44	0.50
1:C:152:MET:HG3	1:C:155:ARG:HH21	1.77	0.50
1:E:156:GLU:HA	1:E:156:GLU:OE1	2.12	0.50
1:B:224:VAL:HG12	1:B:225:GLU:N	2.27	0.49
1:D:224:VAL:O	1:D:228:GLN:HG2	2.12	0.49
1:C:174:GLU:CD	1:D:175:ARG:HH11	2.19	0.49
1:E:160:GLN:HG3	1:F:161:VAL:HG21	1.95	0.49
1:D:168:LYS:O	1:D:172:GLU:HG3	2.13	0.48
1:H:159:ARG:CZ	1:H:159:ARG:HB3	2.42	0.48
1:A:171:VAL:HG12	1:B:171:VAL:CG1	2.32	0.48
1:E:163:GLN:HG2	1:E:167:ASP:OD2	2.14	0.48
1:F:217:ARG:HB2	1:F:217:ARG:HH11	1.75	0.47
1:A:196:ARG:CZ	1:B:192:GLU:OE2	2.63	0.47
1:E:160:GLN:HG3	1:F:161:VAL:CG2	2.44	0.47
1:A:193:MET:HG2	1:B:192:GLU:CD	2.40	0.46
1:G:168:LYS:HE2	1:H:167:ASP:HB3	1.98	0.46
1:B:214:SER:HA	1:B:217:ARG:NH1	2.31	0.46
1:E:217:ARG:O	1:E:221:GLU:HG3	2.16	0.45
1:B:217:ARG:HG2	1:B:221:GLU:OE2	2.15	0.45
1:G:192:GLU:HG2	1:H:193:MET:HG2	1.99	0.45
1:B:204:GLN:O	1:B:208:GLN:HG3	2.16	0.44
1:D:178:LEU:O	1:D:182:ILE:HG13	2.18	0.44
1:G:157:LEU:HB3	1:H:157:LEU:HD13	1.99	0.44
1:A:157:LEU:HD11	1:B:158:ARG:NH2	2.33	0.44
1:G:168:LYS:NZ	1:H:164:LEU:HD23	2.33	0.44
1:H:153:GLU:N	3:H:302:HOH:O	2.50	0.43
1:E:196:ARG:HD3	1:F:192:GLU:OE2	2.17	0.43
1:E:190:GLN:HA	1:E:193:MET:CE	2.43	0.43
1:B:175:ARG:NH2	1:B:176:ASP:OD1	2.48	0.43
1:B:185:LEU:HA	1:B:185:LEU:HD23	1.80	0.43
1:D:217:ARG:HB2	1:D:217:ARG:NH2	2.34	0.43
1:G:192:GLU:HG2	1:H:193:MET:CG	2.49	0.42
1:A:154:MET:HE1	1:B:158:ARG:HH22	1.85	0.42
1:A:175:ARG:HG2	1:B:178:LEU:HD12	2.02	0.42
1:G:222:ARG:HA	1:G:222:ARG:HD2	1.87	0.42
1:A:236:LYS:HA	1:A:236:LYS:HD3	1.87	0.42
1:G:168:LYS:HE2	1:H:167:ASP:CB	2.50	0.42
1:G:220:LEU:CD1	1:H:217:ARG:HG3	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:236:LYS:O	1:D:237:LEU:HB2	2.20	0.41
1:B:225:GLU:O	1:B:226:SER:C	2.61	0.41
1:F:153:GLU:O	1:F:157:LEU:HG	2.20	0.41
1:C:175:ARG:HD2	1:D:174:GLU:OE2	2.20	0.41
1:A:222:ARG:HA	1:A:222:ARG:HD2	1.94	0.41
1:C:217:ARG:O	1:C:221:GLU:HG3	2.21	0.41
1:A:179:ALA:O	1:A:182:ILE:HG22	2.21	0.41
1:H:182:ILE:O	1:H:186:ARG:HG3	2.20	0.40
1:B:186:ARG:O	1:B:190:GLN:HG2	2.21	0.40
1:E:178:LEU:HD12	1:F:175:ARG:HG2	2.03	0.40
1:E:192:GLU:HG2	1:F:193:MET:CG	2.50	0.40
1:H:155:ARG:HB3	1:H:155:ARG:NH1	2.36	0.40
1:C:220:LEU:HD12	1:D:217:ARG:HG3	1.97	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	84/90 (93%)	83 (99%)	1 (1%)	0	100	100
1	B	80/90 (89%)	79 (99%)	1 (1%)	0	100	100
1	C	86/90 (96%)	86 (100%)	0	0	100	100
1	D	83/90 (92%)	82 (99%)	1 (1%)	0	100	100
1	E	84/90 (93%)	83 (99%)	1 (1%)	0	100	100
1	F	82/90 (91%)	82 (100%)	0	0	100	100
1	G	86/90 (96%)	85 (99%)	1 (1%)	0	100	100
1	H	82/90 (91%)	82 (100%)	0	0	100	100
All	All	667/720 (93%)	662 (99%)	5 (1%)	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	65/83 (78%)	65 (100%)	0	100	100
1	B	65/83 (78%)	65 (100%)	0	100	100
1	C	70/83 (84%)	70 (100%)	0	100	100
1	D	70/83 (84%)	70 (100%)	0	100	100
1	E	60/83 (72%)	60 (100%)	0	100	100
1	F	62/83 (75%)	62 (100%)	0	100	100
1	G	70/83 (84%)	70 (100%)	0	100	100
1	H	69/83 (83%)	69 (100%)	0	100	100
All	All	531/664 (80%)	531 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	212	ASN
1	B	166	ASN
1	B	212	ASN
1	C	160	GLN
1	F	166	ASN
1	F	204	GLN
1	H	166	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand 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	$\# Z > 2$
2	GOL	G	301	-	5,5,5	0.66	0	5,5,5	1.03	0

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
2	GOL	G	301	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	301	GOL	O1-C1-C2-C3
2	G	301	GOL	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	86/90 (95%)	0.71	6 (6%)	22 21	24, 36, 57, 66	0
1	B	82/90 (91%)	0.81	8 (9%)	13 12	25, 39, 54, 65	0
1	C	88/90 (97%)	0.50	2 (2%)	61 61	23, 33, 49, 60	0
1	D	85/90 (94%)	0.47	2 (2%)	59 61	20, 34, 50, 60	0
1	E	86/90 (95%)	0.90	8 (9%)	14 13	24, 39, 55, 67	0
1	F	84/90 (93%)	0.71	2 (2%)	59 61	22, 39, 51, 67	0
1	G	88/90 (97%)	0.51	5 (5%)	29 27	20, 33, 51, 64	0
1	H	84/90 (93%)	0.48	2 (2%)	59 61	20, 34, 47, 62	0
All	All	683/720 (94%)	0.63	35 (5%)	33 32	20, 36, 54, 67	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	154	MET	5.6
1	D	237	LEU	5.0
1	A	219	ASP	4.3
1	A	237	LEU	4.0
1	E	237	LEU	3.9
1	C	149	GLY	3.3
1	H	153	GLU	3.3
1	E	152	MET	2.9
1	E	176	ASP	2.9
1	G	214	SER	2.8
1	G	237	LEU	2.8
1	E	154	MET	2.7
1	B	206	PHE	2.6
1	F	152	MET	2.6
1	E	158	ARG	2.4
1	A	155	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	179	ALA	2.4
1	E	233	PHE	2.4
1	B	194	LEU	2.4
1	A	165	THR	2.3
1	H	206	PHE	2.3
1	G	151	HIS	2.3
1	G	176	ASP	2.3
1	B	213	ALA	2.3
1	B	190	GLN	2.3
1	A	174	GLU	2.3
1	E	202	THR	2.2
1	G	150	PRO	2.2
1	B	158	ARG	2.1
1	A	163	GLN	2.1
1	F	219	ASP	2.1
1	B	233	PHE	2.1
1	E	206	PHE	2.1
1	C	214	SER	2.1
1	D	155	ARG	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

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
2	GOL	G	301	6/6	0.77	0.13	45,55,65,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.