



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:16 PM UTC

PDB ID : 8GJF / pdb_00008gjf
Title : afupcna bound with peptide mimetic
Authors : Vandborg, B.; Bruning, J.B.
Deposited on : 2023-03-15
Resolution : 2.00 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
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

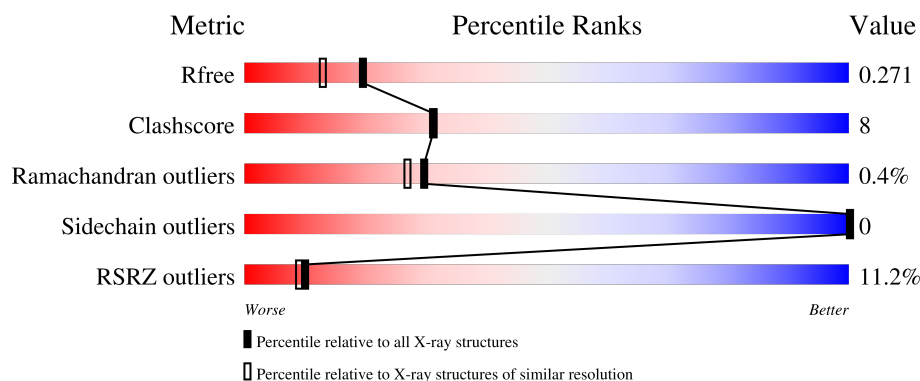
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	256	12% 85% 14%
1	B	256	10% 84% 15%
1	C	256	12% 80% 20%
2	D	15	87% 13%
2	E	15	13% 60% 13% 7% 20%

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Mol	Chain	Length	Quality of chain
2	F	15	 7% 80% 13% 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6392 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 Proliferating cell nuclear antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	1	0
			1867	1179	311	365	12			
1	B	254	Total	C	N	O	S	0	1	0
			1885	1187	320	365	13			
1	C	255	Total	C	N	O	S	0	0	0
			1892	1195	309	375	13			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	59	ALA	PRO	conflict	UNP A0A229Y5V5
B	59	ALA	PRO	conflict	UNP A0A229Y5V5
C	59	ALA	PRO	conflict	UNP A0A229Y5V5

- Molecule 2 is a protein called LYS-ARG-ARG-GLN-THR-SER-MET-THR-ASP-PHE-TYR-HIS-SER-LYS-ARG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	13	Total	C	N	O	S	0	0	0
			111	67	22	21	1			
2	E	12	Total	C	N	O	S	0	0	0
			100	61	18	20	1			
2	F	14	Total	C	N	O	S	0	0	0
			116	72	21	22	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	129	Total	O	0	0
			129	129		
3	B	123	Total	O	0	0
			123	123		

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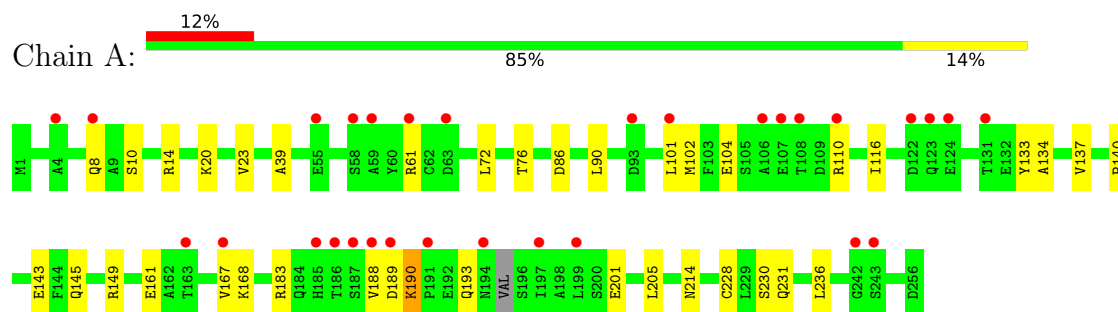
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	131	Total 131	O 131	0	0
3	D	16	Total 16	O 16	0	0
3	E	6	Total 6	O 6	0	0
3	F	16	Total 16	O 16	0	0

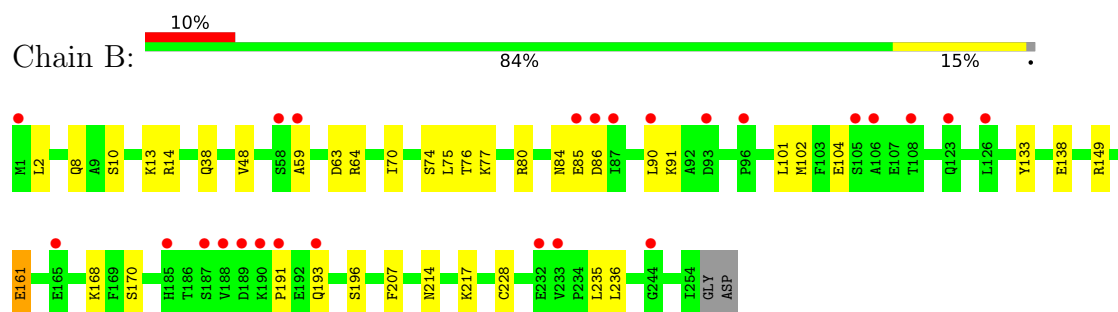
3 Residue-property plots [i](#)

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.

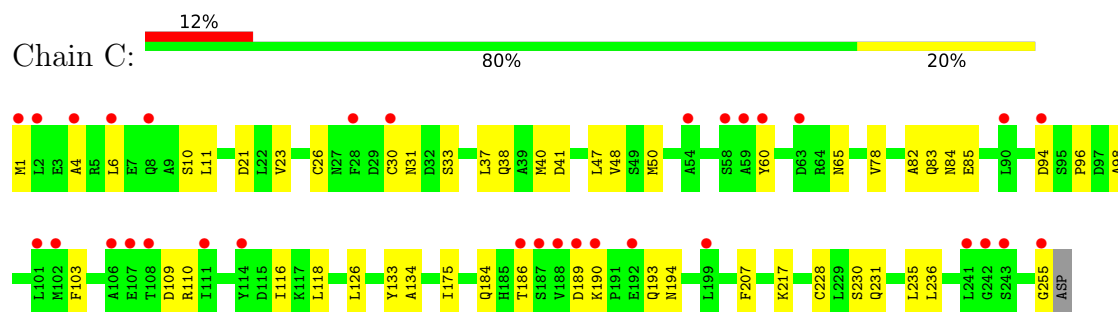
- Molecule 1: Proliferating cell nuclear antigen



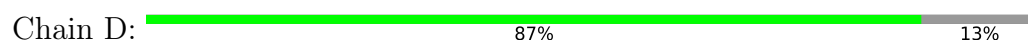
- Molecule 1: Proliferating cell nuclear antigen



- Molecule 1: Proliferating cell nuclear antigen



- Molecule 2: LYS-ARG-ARG-GLN-THR-SER-MET-THR-ASP-PHE-TYR-HIS-SER-LYS-ARG

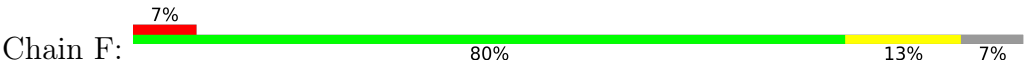




● Molecule 2: LYS-ARG-ARG-GLN-THR-SER-MET-THR-ASP-PHE-TYR-HIS-SER-LYS-ARG



● Molecule 2: LYS-ARG-ARG-GLN-THR-SER-MET-THR-ASP-PHE-TYR-HIS-SER-LYS-ARG



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	146.00Å 84.83Å 70.45Å 90.00° 91.41° 90.00°	Depositor
Resolution (Å)	42.20 – 2.00 42.20 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.1 (42.20-2.00) 99.1 (42.20-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.233 , 0.270 0.234 , 0.271	Depositor DCC
R_{free} test set	2889 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	38.3	Xtriage
Anisotropy	0.286	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.000 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.085 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.053 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.015 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6392	wwPDB-VP
Average B, all atoms (Å ²)	49.0	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

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	0.22	0/1892	0.79	3/2571 (0.1%)
1	B	0.32	0/1911	0.58	1/2595 (0.0%)
1	C	0.23	0/1916	0.51	1/2600 (0.0%)
2	D	0.14	0/113	0.44	0/150
2	E	3.81	1/102 (1.0%)	0.43	0/136
2	F	0.08	0/118	0.29	0/156
All	All	0.56	1/6052 (0.0%)	0.63	5/8208 (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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	153	SER	C-N	-38.45	0.79	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	GLU	OE1-CD-OE2	-18.93	77.46	122.90
1	A	201	GLU	CG-CD-OE1	18.31	160.52	118.40
1	A	201	GLU	CG-CD-OE2	-15.36	83.06	118.40
1	B	161	GLU	CG-CD-OE2	-6.09	104.39	118.40
1	C	184	GLN	CA-CB-CG	5.17	124.43	114.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	161	GLU	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	1867	0	1815	31	1
1	B	1885	0	1863	24	0
1	C	1892	0	1868	39	1
2	D	111	0	97	0	0
2	E	100	0	83	7	0
2	F	116	0	103	2	0
3	A	129	0	0	5	0
3	B	123	0	0	3	0
3	C	131	0	0	9	0
3	D	16	0	0	0	0
3	E	6	0	0	1	0
3	F	16	0	0	0	0
All	All	6392	0	5829	97	1

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 8.

All (97) 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)
2:E:153:SER:C	2:E:154:LYS:CA	1.97	1.34
2:E:153:SER:O	2:E:154:LYS:N	1.75	1.19
2:E:153:SER:CA	2:E:154:LYS:N	2.05	1.18
2:E:153:SER:O	2:E:154:LYS:CA	2.10	0.92
2:E:153:SER:O	2:E:154:LYS:HA	1.70	0.91
2:E:153:SER:C	2:E:154:LYS:N	0.79	0.89
1:C:26:CYS:SG	3:C:309:HOH:O	2.40	0.79
1:A:20:LYS:HD3	1:A:76:THR:HG21	1.71	0.73
1:B:86:ASP:N	3:B:301:HOH:O	2.27	0.66
1:B:76:THR:O	1:B:80:ARG:HG3	1.98	0.64
1:A:161:GLU:HG3	1:A:168:LYS:HB3	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:ARG:NH2	3:A:302:HOH:O	2.32	0.62
1:C:1:MET:HA	1:C:1:MET:HE2	1.81	0.62
1:A:145:GLN:NE2	3:A:302:HOH:O	2.32	0.62
1:B:63:ASP:O	1:B:64:ARG:HD3	2.01	0.61
1:C:60:TYR:OH	3:C:301:HOH:O	2.16	0.60
1:B:14:ARG:NH1	3:B:303:HOH:O	2.33	0.60
1:A:8:GLN:N	1:A:8:GLN:OE1	2.34	0.60
1:B:8:GLN:NE2	1:B:85[A]:GLU:OE1	2.33	0.60
1:C:231:GLN:NE2	3:C:306:HOH:O	2.33	0.59
1:A:134:ALA:H	1:A:230:SER:HB3	1.67	0.59
1:A:8:GLN:HG2	1:A:8:GLN:O	2.05	0.56
1:C:31:ASN:O	1:C:60:TYR:OH	2.20	0.56
1:C:4:ALA:HA	3:C:357:HOH:O	2.06	0.56
1:A:110[B]:ARG:NH2	3:A:306:HOH:O	2.36	0.56
1:C:10:SER:HB3	1:C:84:ASN:HB3	1.88	0.56
1:A:205:LEU:HD21	1:A:231:GLN:HA	1.88	0.55
1:A:86:ASP:OD2	1:A:110[A]:ARG:NH1	2.40	0.55
1:B:228:CYS:HB2	1:B:236:LEU:HB3	1.90	0.54
1:B:138:GLU:HB2	1:B:196:SER:HB3	1.91	0.53
1:C:186:THR:HA	1:C:194:ASN:OD1	2.08	0.52
1:C:6:LEU:HD11	1:C:11:LEU:HD23	1.91	0.52
2:E:143:ARG:N	3:E:202:HOH:O	2.43	0.51
1:C:30:CYS:HB3	3:C:301:HOH:O	2.08	0.51
1:C:65:ASN:ND2	3:C:303:HOH:O	2.28	0.51
1:C:255:GLY:HA3	2:F:142:ARG:HB2	1.94	0.50
1:A:137:VAL:HG21	1:A:167:VAL:HG21	1.94	0.50
1:A:228:CYS:HB2	1:A:236:LEU:HB3	1.94	0.50
1:C:38:GLN:HA	1:C:48:VAL:O	2.12	0.49
1:A:10:SER:O	1:A:14:ARG:HG3	2.13	0.48
1:A:145:GLN:O	1:A:149:ARG:HG3	2.13	0.48
1:C:78:VAL:HG21	1:C:116:ILE:CD1	2.44	0.48
1:A:183:ARG:HB2	1:C:109:ASP:HB3	1.96	0.47
1:A:188:VAL:O	1:A:190:LYS:N	2.48	0.47
1:A:102:MET:HE3	1:A:104:GLU:CD	2.38	0.47
1:B:38:GLN:HA	1:B:48:VAL:O	2.14	0.47
1:C:40:MET:HG3	1:C:47:LEU:HD12	1.97	0.47
1:B:214:ASN:OD1	1:B:217:LYS:NZ	2.46	0.47
1:A:110[B]:ARG:HE	1:A:110[B]:ARG:HB3	1.64	0.47
1:B:133:TYR:CG	1:B:228:CYS:HB3	2.50	0.47
1:C:85:GLU:OE1	1:C:85:GLU:N	2.42	0.46
1:C:126:LEU:O	2:F:154:LYS:HE2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:134:ALA:H	1:C:230:SER:HB3	1.79	0.46
1:B:13:LYS:HE2	1:B:84:ASN:HD22	1.79	0.46
1:A:90:LEU:CD2	1:A:101:LEU:HD23	2.46	0.46
1:B:207:PHE:CZ	1:B:235:LEU:HB2	2.51	0.46
1:A:140:PRO:HD3	1:A:193:GLN:O	2.15	0.46
1:A:214:ASN:ND2	3:A:313:HOH:O	2.48	0.46
1:C:83:GLN:HB3	1:C:85:GLU:OE1	2.16	0.46
1:B:13:LYS:HE2	1:B:84:ASN:ND2	2.31	0.45
1:B:74:SER:HB3	1:C:175:ILE:HG23	1.99	0.45
1:A:72:LEU:O	1:A:76:THR:HG23	2.17	0.45
1:C:50:MET:HE2	1:C:50:MET:HB3	1.92	0.45
1:C:82:ALA:HB2	1:C:103:PHE:CD2	2.52	0.45
1:C:207:PHE:CZ	1:C:235:LEU:HB2	2.52	0.45
1:C:98:ALA:HA	1:C:118:LEU:HG	1.99	0.44
1:C:193:GLN:N	3:C:318:HOH:O	2.50	0.44
1:A:143:GLU:OE1	1:C:110:ARG:NH2	2.44	0.44
1:A:8:GLN:H	1:A:8:GLN:CD	2.26	0.44
1:C:228:CYS:HB2	1:C:236:LEU:HB3	1.98	0.44
1:A:133:TYR:CG	1:A:228:CYS:HB3	2.53	0.44
1:A:23:VAL:HG11	1:A:39:ALA:HB1	1.99	0.43
1:C:10:SER:HB3	1:C:84:ASN:CB	2.47	0.43
1:B:90:LEU:HD22	1:B:101:LEU:HD23	2.00	0.43
1:B:168:LYS:HE2	1:B:170:SER:HB3	2.00	0.43
1:A:161:GLU:HG2	1:A:168:LYS:O	2.19	0.43
1:B:10:SER:O	1:B:14:ARG:HD3	2.19	0.43
1:C:94:ASP:C	1:C:96:PRO:HD3	2.43	0.43
1:B:2:LEU:O	1:B:91:LYS:HA	2.18	0.43
1:B:191:PRO:C	1:B:193:GLN:H	2.27	0.42
1:C:21:ASP:OD2	1:C:217:LYS:NZ	2.48	0.42
1:C:1:MET:HG2	1:C:94:ASP:OD1	2.20	0.42
1:C:30:CYS:SG	3:C:414:HOH:O	2.62	0.42
1:C:189:ASP:O	1:C:190:LYS:HE2	2.20	0.42
1:A:101:LEU:HD11	1:A:116:ILE:HD11	2.01	0.42
1:C:23:VAL:HG12	1:C:41:ASP:HA	2.00	0.42
1:A:161:GLU:CG	1:A:168:LYS:HB3	2.50	0.41
1:A:183:ARG:HD3	1:A:183:ARG:HA	1.77	0.41
1:C:133:TYR:CG	1:C:228:CYS:HB3	2.56	0.41
1:B:77:LYS:HD2	1:C:175:ILE:CD1	2.51	0.41
1:C:37:LEU:HB3	1:C:50:MET:HE2	2.03	0.41
1:B:102:MET:HE3	1:B:104:GLU:CD	2.46	0.41
1:C:60:TYR:HD1	3:C:357:HOH:O	2.05	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:ARG:NH2	3:B:311:HOH:O	2.54	0.40
1:A:20:LYS:NZ	3:A:312:HOH:O	2.47	0.40
1:B:70:ILE:HG21	1:B:75:LEU:HD22	2.03	0.40

All (1) 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:61:ARG:NH1	1:C:33:SER:OG[1_545]	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	254/256 (99%)	248 (98%)	4 (2%)	2 (1%)	16	11
1	B	253/256 (99%)	244 (96%)	8 (3%)	1 (0%)	30	27
1	C	253/256 (99%)	247 (98%)	6 (2%)	0	100	100
2	D	11/15 (73%)	11 (100%)	0	0	100	100
2	E	10/15 (67%)	10 (100%)	0	0	100	100
2	F	12/15 (80%)	12 (100%)	0	0	100	100
All	All	793/813 (98%)	772 (97%)	18 (2%)	3 (0%)	30	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	ASP
1	A	190	LYS
1	B	59	ALA

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	192/221 (87%)	192 (100%)	0	100	100
1	B	201/221 (91%)	201 (100%)	0	100	100
1	C	203/221 (92%)	203 (100%)	0	100	100
2	D	12/15 (80%)	12 (100%)	0	100	100
2	E	11/15 (73%)	11 (100%)	0	100	100
2	F	12/15 (80%)	12 (100%)	0	100	100
All	All	631/708 (89%)	631 (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 (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	84	ASN
1	A	152	ASN
1	A	184	GLN
1	B	8	GLN
1	B	24	GLN
1	B	27	ASN
1	B	84	ASN
1	B	145	GLN
1	C	24	GLN
1	C	27	ASN
1	C	125	HIS
2	D	152	HIS

5.3.3 RNA [i](#)

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](#)

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	E	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	153:SER	C	154:LYS	N	0.79

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	255/256 (99%)	0.73	30 (11%) 9 8	27, 44, 82, 123	1 (0%)
1	B	254/256 (99%)	0.79	25 (9%) 13 12	33, 45, 89, 119	1 (0%)
1	C	255/256 (99%)	0.94	32 (12%) 8 7	27, 45, 85, 131	0
2	D	13/15 (86%)	0.57	0 100 100	30, 37, 62, 74	0
2	E	12/15 (80%)	0.97	2 (16%) 4 4	44, 53, 66, 73	0
2	F	14/15 (93%)	0.66	1 (7%) 22 20	32, 38, 67, 69	0
All	All	803/813 (98%)	0.82	90 (11%) 10 9	27, 45, 83, 131	2 (0%)

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	58	SER	6.0
1	A	188	VAL	4.8
1	C	4	ALA	4.8
1	C	188	VAL	4.7
1	C	59	ALA	4.2
1	B	187	SER	4.1
1	C	1	MET	4.1
1	C	255	GLY	3.9
1	C	63	ASP	3.9
1	A	186	THR	3.9
1	B	188	VAL	3.8
1	B	106	ALA	3.7
1	A	59	ALA	3.6
1	C	60	TYR	3.6
1	B	189	ASP	3.5
1	B	59	ALA	3.5
1	C	2	LEU	3.4
1	B	193	GLN	3.4
1	C	90	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	189	ASP	3.2
1	A	187	SER	3.2
1	A	58	SER	3.2
2	F	142	ARG	3.1
1	C	241	LEU	3.1
1	C	187	SER	3.0
1	A	197	ILE	3.0
1	A	107	GLU	3.0
1	C	190	LYS	3.0
1	C	102	MET	2.9
1	A	106	ALA	2.9
1	C	54	ALA	2.9
1	C	6	LEU	2.8
1	B	86	ASP	2.8
1	A	189	ASP	2.8
1	B	190	LYS	2.8
1	A	243	SER	2.8
1	B	232	GLU	2.8
1	C	108	THR	2.8
1	A	123	GLN	2.7
1	A	63	ASP	2.7
1	C	186	THR	2.7
1	A	163	THR	2.7
1	C	94	ASP	2.6
1	A	4	ALA	2.6
1	B	185	HIS	2.6
1	C	111	ILE	2.6
1	A	122	ASP	2.5
1	C	106	ALA	2.5
1	A	93	ASP	2.4
1	B	1	MET	2.4
1	B	165	GLU	2.4
1	C	101	LEU	2.4
1	C	242	GLY	2.4
1	C	30	CYS	2.4
1	B	108	THR	2.4
1	B	96	PRO	2.4
1	A	110[A]	ARG	2.4
1	A	131	THR	2.4
1	A	8	GLN	2.3
1	A	191	PRO	2.3
2	E	147	MET	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	85[A]	GLU	2.3
1	B	191	PRO	2.3
1	B	244	GLY	2.3
2	E	152	HIS	2.3
1	C	8	GLN	2.3
1	A	101	LEU	2.3
1	B	126	LEU	2.3
1	A	185	HIS	2.2
1	C	243	SER	2.2
1	C	114	TYR	2.2
1	B	58	SER	2.2
1	A	61	ARG	2.2
1	B	105	SER	2.2
1	A	124	GLU	2.1
1	B	87	ILE	2.1
1	C	192	GLU	2.1
1	A	55	GLU	2.1
1	B	93	ASP	2.1
1	C	107	GLU	2.1
1	A	167	VAL	2.1
1	A	242	GLY	2.1
1	A	199	LEU	2.1
1	B	123	GLN	2.1
1	C	28	PHE	2.1
1	B	90	LEU	2.0
1	A	108	THR	2.0
1	B	233	VAL	2.0
1	A	194	ASN	2.0
1	C	199	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.