



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:31 PM UTC

PDB ID : 6N2G / pdb_00006n2g
Title : Crystal structure of Caenorhabditis elegans NAP1
Authors : Bhattacharyya, S.; DArcy, S.
Deposited on : 2018-11-13
Resolution : 3.00 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

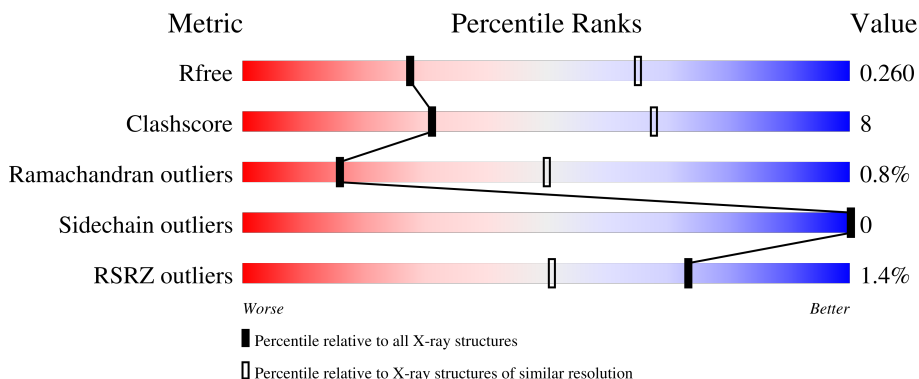
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	316	% 73% 17% 10%
1	B	316	2% 65% 22% 14%
1	C	316	% 70% 19% 11%
1	D	316	% 72% 14% 14%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 8866 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.

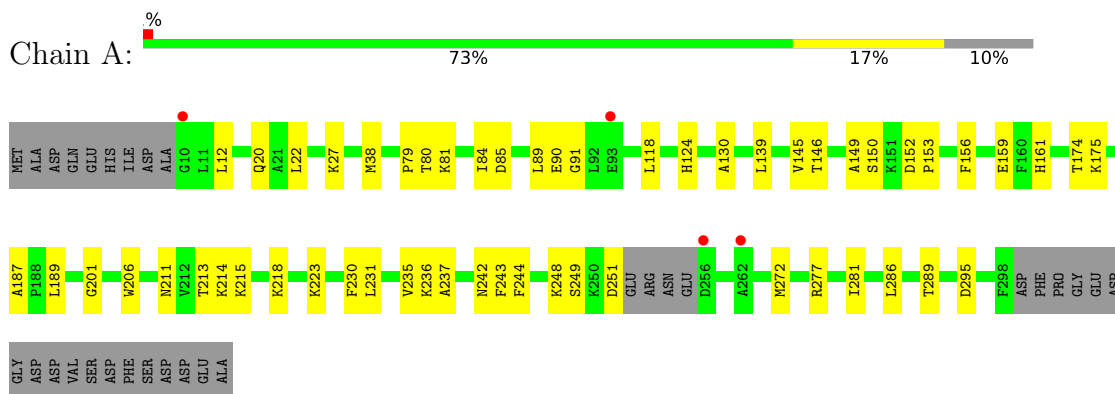
- Molecule 1 is a protein called Nucleosome Assembly Protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2268	1453	367	443	5			
1	B	273	Total	C	N	O	S	0	0	0
			2180	1401	355	420	4			
1	C	281	Total	C	N	O	S	0	0	0
			2241	1440	363	433	5			
1	D	273	Total	C	N	O	S	0	0	0
			2177	1400	355	418	4			

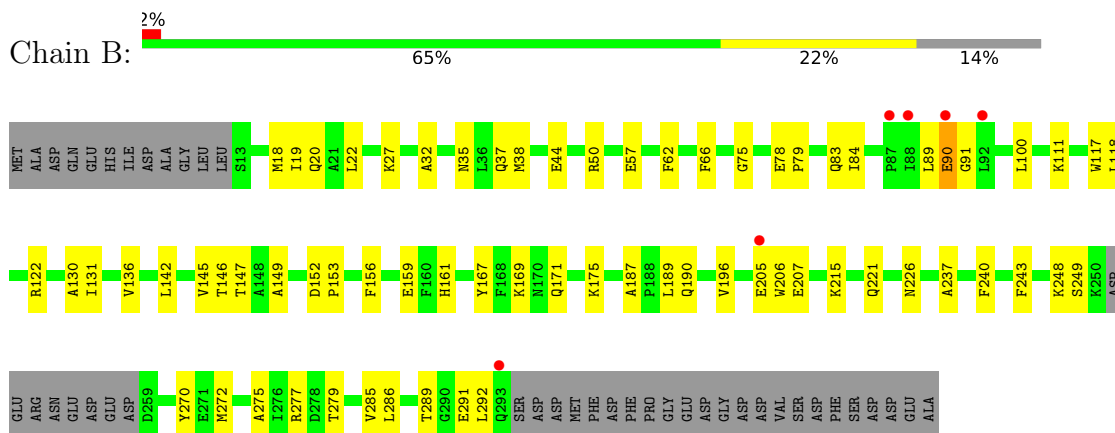
3 Residue-property plots

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.

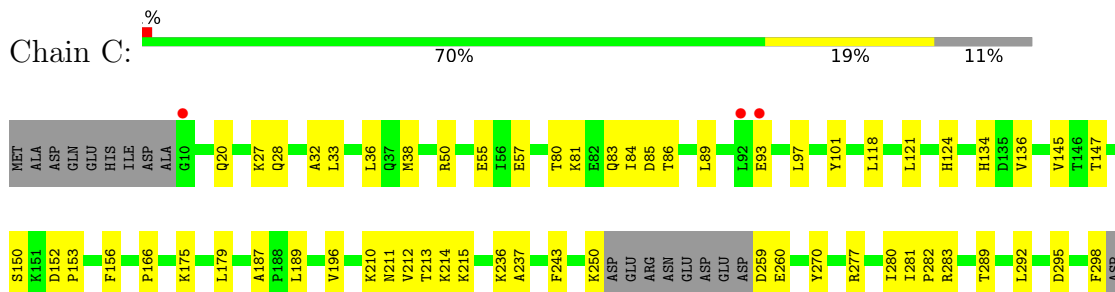
• Molecule 1: Nucleosome Assembly Protein



• Molecule 1: Nucleosome Assembly Protein



• Molecule 1: Nucleosome Assembly Protein



PHE
PRO
GLY
GLY
ASP
GLY
ASP
ASP
VAL
SER
ASP
PHE
SER
ASP
ASP
GLU
ALA

● Molecule 1: Nucleosome Assembly Protein



PHE
SER
ASP
ASP
GLU
ALA

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	125.60Å 154.87Å 92.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.80 – 3.00 19.80 – 3.00	Depositor EDS
% Data completeness (in resolution range)	88.6 (19.80-3.00) 89.3 (19.80-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.98Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.241 , 0.255 0.247 , 0.260	Depositor DCC
R_{free} test set	2016 reflections (6.14%)	wwPDB-VP
Wilson B-factor (Å ²)	58.3	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 28.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8866	wwPDB-VP
Average B, all atoms (Å ²)	56.0	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 40.14 % 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 2.8608e-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 [i](#)

5.1 Standard geometry [i](#)

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	0.17	0/2315	0.40	0/3126
1	B	0.17	0/2226	0.43	0/3006
1	C	0.18	0/2288	0.42	0/3089
1	D	0.16	0/2223	0.45	0/3002
All	All	0.17	0/9052	0.42	0/12223

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2268	0	2225	35	0
1	B	2180	0	2162	58	0
1	C	2241	0	2218	48	0
1	D	2177	0	2160	26	0
All	All	8866	0	8765	144	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 8.

All (144) 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:D:18:MET:O	1:D:22:LEU:HD13	1.72	0.90
1:B:18:MET:O	1:B:22:LEU:HD13	1.76	0.85
1:B:152:ASP:OD1	1:B:153:PRO:CD	2.25	0.84
1:B:152:ASP:CG	1:B:153:PRO:HD3	2.05	0.81
1:B:152:ASP:OD1	1:B:153:PRO:HD2	1.82	0.80
1:A:215:LYS:HB2	1:A:237:ALA:HB2	1.67	0.76
1:D:215:LYS:HB2	1:D:237:ALA:HB2	1.68	0.76
1:A:150:SER:OG	1:C:250:LYS:NZ	2.22	0.72
1:B:50:ARG:NH1	1:C:57:GLU:OE2	2.23	0.71
1:D:81:LYS:NZ	1:D:85:ASP:OD1	2.24	0.69
1:D:213:THR:OG1	1:D:242:ASN:OD1	2.10	0.68
1:A:124:HIS:NE2	1:A:295:ASP:OD1	2.25	0.67
1:C:215:LYS:HB2	1:C:237:ALA:HB2	1.76	0.67
1:C:213:THR:HB	1:C:236:LYS:HE3	1.76	0.67
1:B:35:ASN:HB2	1:C:86:THR:HG21	1.78	0.65
1:B:75:GLY:O	1:B:111:LYS:NZ	2.30	0.64
1:D:147:THR:OG1	1:D:277:ARG:NH1	2.30	0.64
1:C:93:GLU:HB3	1:C:97:LEU:HB2	1.79	0.63
1:B:152:ASP:OD1	1:B:153:PRO:HD3	1.95	0.63
1:A:146:THR:HG22	1:A:159:GLU:HB2	1.80	0.62
1:A:159:GLU:HG2	1:A:174:THR:HG23	1.81	0.62
1:C:211:ASN:OD1	1:C:213:THR:OG1	2.16	0.61
1:C:156:PHE:CD1	1:C:277:ARG:HD2	2.35	0.61
1:B:275:ALA:O	1:B:279:THR:OG1	2.17	0.61
1:A:20:GLN:O	1:A:27:LYS:NZ	2.31	0.61
1:B:152:ASP:CG	1:B:153:PRO:CD	2.72	0.61
1:B:79:PRO:HB3	1:C:32:ALA:HB2	1.82	0.60
1:D:211:ASN:OD1	1:D:213:THR:HG22	2.01	0.60
1:B:83:GLN:HB3	1:C:32:ALA:HA	1.82	0.60
1:B:285:VAL:HG21	1:C:33:LEU:HB3	1.83	0.60
1:B:146:THR:HG22	1:B:159:GLU:HB2	1.84	0.59
1:C:55:GLU:OE1	1:C:277:ARG:NH1	2.32	0.59
1:B:44:GLU:OE1	1:C:283:ARG:NH2	2.32	0.57
1:A:211:ASN:OD1	1:A:213:THR:HG22	2.04	0.57
1:C:145:VAL:HG12	1:C:281:ILE:HD11	1.85	0.57
1:B:90:GLU:N	1:B:291:GLU:OE2	2.38	0.56
1:B:84:ILE:HD12	1:C:28:GLN:HB3	1.87	0.56
1:B:248:LYS:HD2	1:B:249:SER:H	1.70	0.56
1:B:122:ARG:HH11	1:B:131:ILE:HG13	1.69	0.56
1:D:208:ASP:O	1:D:210:LYS:N	2.38	0.56
1:B:147:THR:HG21	1:B:277:ARG:HG3	1.87	0.55
1:B:286:LEU:HA	1:B:289:THR:HG22	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:GLY:O	1:D:171:GLN:NE2	2.38	0.54
1:A:80:THR:O	1:A:84:ILE:HG12	2.07	0.54
1:C:150:SER:OG	1:C:153:PRO:HD2	2.07	0.54
1:B:156:PHE:CD1	1:B:277:ARG:HD2	2.43	0.54
1:B:190:GLN:OE1	1:B:190:GLN:N	2.41	0.54
1:B:38:MET:HG3	1:B:189:LEU:HD21	1.90	0.53
1:B:57:GLU:OE2	1:C:50:ARG:NH2	2.40	0.53
1:B:190:GLN:HA	1:C:298:PHE:CE2	2.44	0.53
1:B:215:LYS:HB2	1:B:237:ALA:HB2	1.91	0.53
1:A:213:THR:OG1	1:A:242:ASN:OD1	2.26	0.52
1:C:97:LEU:HD21	1:C:101:TYR:CE2	2.44	0.52
1:A:81:LYS:NZ	1:A:85:ASP:OD2	2.40	0.52
1:A:215:LYS:N	1:A:235:VAL:O	2.42	0.52
1:D:275:ALA:O	1:D:279:THR:OG1	2.24	0.52
1:C:121:LEU:HD23	1:C:280:ILE:HD13	1.93	0.51
1:A:145:VAL:HG12	1:A:281:ILE:HD11	1.93	0.51
1:C:166:PRO:O	1:C:210:LYS:HD2	2.11	0.51
1:A:175:LYS:NZ	1:A:243:PHE:O	2.37	0.50
1:C:175:LYS:NZ	1:C:243:PHE:O	2.35	0.50
1:A:156:PHE:CD1	1:A:277:ARG:HD2	2.47	0.50
1:C:32:ALA:O	1:C:36:LEU:HD13	2.12	0.49
1:B:79:PRO:HB2	1:B:84:ILE:HD11	1.94	0.49
1:B:131:ILE:HG22	1:B:240:PHE:HB2	1.94	0.49
1:B:122:ARG:HA	1:B:131:ILE:HD11	1.95	0.49
1:B:130:ALA:HB2	1:B:272:MET:HE1	1.95	0.48
1:D:187:ALA:HB1	1:D:190:GLN:HG3	1.95	0.48
1:B:19:ILE:HA	1:B:22:LEU:HD22	1.94	0.48
1:B:100:LEU:O	1:C:27:LYS:HE3	2.14	0.48
1:B:286:LEU:HB3	1:B:292:LEU:HG	1.96	0.48
1:C:147:THR:HG21	1:C:277:ARG:HG3	1.95	0.48
1:B:117:TRP:CZ3	1:B:145:VAL:HG22	2.48	0.48
1:A:248:LYS:NZ	1:A:251:ASP:HB2	2.29	0.47
1:B:175:LYS:HG2	1:B:243:PHE:CE2	2.49	0.47
1:B:78:GLU:HG3	1:B:79:PRO:HD2	1.96	0.47
1:D:135:ASP:OD2	1:D:240:PHE:N	2.37	0.47
1:A:206:TRP:CE3	1:A:211:ASN:HA	2.49	0.47
1:D:262:ALA:HA	1:D:265:PHE:HD2	1.78	0.47
1:A:152:ASP:HB3	1:A:153:PRO:HD3	1.97	0.47
1:D:286:LEU:HA	1:D:289:THR:HG22	1.97	0.47
1:C:124:HIS:NE2	1:C:295:ASP:OD1	2.40	0.47
1:A:89:LEU:C	1:A:91:GLY:H	2.22	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:LYS:HE2	1:A:236:LYS:HZ3	1.80	0.46
1:B:62:PHE:HB3	1:B:66:PHE:CE2	2.50	0.46
1:D:175:LYS:NZ	1:D:243:PHE:O	2.47	0.46
1:A:161:HIS:ND1	1:C:152:ASP:OD1	2.49	0.46
1:D:152:ASP:HB3	1:D:153:PRO:HD3	1.97	0.46
1:D:145:VAL:HG12	1:D:281:ILE:HD11	1.97	0.46
1:B:89:LEU:C	1:B:91:GLY:H	2.23	0.46
1:B:32:ALA:HA	1:C:83:GLN:HB3	1.98	0.45
1:B:100:LEU:HD11	1:C:20:GLN:HA	1.97	0.45
1:C:38:MET:HG3	1:C:189:LEU:HD21	1.98	0.45
1:C:80:THR:O	1:C:84:ILE:HG12	2.16	0.45
1:D:160:PHE:O	1:D:172:VAL:HA	2.17	0.45
1:A:230:PHE:H	1:D:226:ASN:HD21	1.65	0.45
1:D:156:PHE:CD1	1:D:277:ARG:HD2	2.52	0.45
1:A:79:PRO:HB2	1:A:84:ILE:HD11	1.99	0.45
1:A:286:LEU:HA	1:A:289:THR:HG22	1.98	0.45
1:D:221:GLN:HG3	1:D:231:LEU:HD23	1.99	0.45
1:B:37:GLN:NE2	1:C:282:PRO:O	2.42	0.45
1:C:214:LYS:HD3	1:C:236:LYS:HD2	1.98	0.45
1:D:18:MET:O	1:D:22:LEU:CD1	2.54	0.45
1:A:235:VAL:HG22	1:A:236:LYS:H	1.80	0.44
1:C:20:GLN:O	1:C:27:LYS:NZ	2.47	0.44
1:A:38:MET:HG3	1:A:189:LEU:HD21	2.00	0.43
1:B:20:GLN:O	1:B:27:LYS:NZ	2.48	0.43
1:C:152:ASP:HB3	1:C:153:PRO:HD3	2.00	0.43
1:A:248:LYS:HD2	1:A:249:SER:H	1.83	0.43
1:C:259:ASP:OD1	1:C:260:GLU:N	2.46	0.43
1:B:90:GLU:HB2	1:C:189:LEU:HD12	2.00	0.43
1:B:118:LEU:HD21	1:B:136:VAL:HG13	2.00	0.42
1:A:12:LEU:HD23	1:A:12:LEU:H	1.83	0.42
1:B:142:LEU:HD12	1:B:161:HIS:O	2.19	0.42
1:A:218:LYS:HA	1:A:231:LEU:O	2.20	0.42
1:C:89:LEU:HD23	1:C:89:LEU:HA	1.84	0.42
1:C:134:HIS:CE1	1:C:212:VAL:HG13	2.54	0.42
1:D:249:SER:O	1:D:250:LYS:HB2	2.18	0.42
1:B:149:ALA:HA	1:B:156:PHE:HA	2.02	0.42
1:A:130:ALA:HB2	1:A:272:MET:HE1	2.02	0.42
1:B:19:ILE:HD11	1:C:289:THR:HA	2.01	0.42
1:B:75:GLY:C	1:B:111:LYS:HZ2	2.28	0.42
1:C:81:LYS:NZ	1:C:85:ASP:OD1	2.52	0.42
1:C:292:LEU:HD21	1:C:298:PHE:CZ	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:179:LEU:HD23	1:C:196:VAL:HA	2.02	0.42
1:B:19:ILE:O	1:B:27:LYS:NZ	2.54	0.41
1:B:169:LYS:HB2	1:B:205:GLU:HB2	2.02	0.41
1:D:89:LEU:HD23	1:D:89:LEU:HA	1.80	0.41
1:B:196:VAL:HG11	1:B:270:TYR:CG	2.55	0.41
1:D:259:ASP:C	1:D:261:GLN:H	2.27	0.41
1:A:22:LEU:HD23	1:A:22:LEU:HA	1.93	0.41
1:A:223:LYS:HD2	1:A:223:LYS:HA	1.91	0.41
1:A:149:ALA:HA	1:A:156:PHE:HA	2.02	0.41
1:D:150:SER:OG	1:D:153:PRO:HD2	2.21	0.41
1:B:167:TYR:HB3	1:B:206:TRP:HZ3	1.85	0.41
1:B:221:GLN:HB2	1:B:226:ASN:O	2.21	0.41
1:C:179:LEU:HD21	1:C:270:TYR:CE1	2.56	0.41
1:A:201:GLY:HA3	1:A:244:PHE:O	2.20	0.40
1:A:152:ASP:HB2	1:C:250:LYS:HD3	2.03	0.40
1:A:118:LEU:HD13	1:A:139:LEU:HB2	2.03	0.40
1:B:285:VAL:HG22	1:C:33:LEU:HD13	2.04	0.40
1:C:118:LEU:HD21	1:C:136:VAL:HG13	2.04	0.40
1:B:90:GLU:HG3	1:C:189:LEU:HG	2.02	0.40
1:B:171:GLN:NE2	1:D:150:SER:HA	2.37	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	281/316 (89%)	273 (97%)	6 (2%)	2 (1%)	18 53
1	B	269/316 (85%)	262 (97%)	4 (2%)	3 (1%)	11 43
1	C	277/316 (88%)	270 (98%)	6 (2%)	1 (0%)	30 65
1	D	269/316 (85%)	261 (97%)	5 (2%)	3 (1%)	11 43
All	All	1096/1264 (87%)	1066 (97%)	21 (2%)	9 (1%)	16 50

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	187	ALA
1	D	209	GLY
1	D	91	GLY
1	A	90	GLU
1	B	90	GLU
1	D	187	ALA
1	B	207	GLU
1	A	187	ALA
1	B	187	ALA

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	242/270 (90%)	242 (100%)	0	100	100
1	B	233/270 (86%)	233 (100%)	0	100	100
1	C	240/270 (89%)	240 (100%)	0	100	100
1	D	232/270 (86%)	232 (100%)	0	100	100
All	All	947/1080 (88%)	947 (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 (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	42	GLN
1	A	52	HIS
1	B	171	GLN
1	C	37	GLN
1	C	96	GLN
1	C	190	GLN
1	D	35	ASN
1	D	221	GLN

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](#)

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	285/316 (90%)	-0.03	4 (1%)	73 51	30, 49, 85, 109	0
1	B	273/316 (86%)	0.38	6 (2%)	62 39	36, 65, 100, 115	0
1	C	281/316 (88%)	-0.00	3 (1%)	78 57	28, 45, 87, 100	0
1	D	273/316 (86%)	0.00	3 (1%)	78 57	35, 53, 83, 103	0
All	All	1112/1264 (87%)	0.09	16 (1%)	73 51	28, 53, 91, 115	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	256	ASP	4.2
1	D	259	ASP	3.6
1	C	92	LEU	3.0
1	B	293	GLN	2.6
1	A	10	GLY	2.5
1	B	92	LEU	2.5
1	D	92	LEU	2.4
1	C	93	GLU	2.3
1	A	93	GLU	2.3
1	D	152	ASP	2.3
1	B	205	GLU	2.2
1	B	88	ILE	2.2
1	A	262	ALA	2.2
1	B	90	GLU	2.1
1	B	87	PRO	2.0
1	C	10	GLY	2.0

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.