



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 06:45 PM UTC

PDB ID : 7JSM / pdb_00007jsm
Title : CRYSTAL STRUCTURE OF NATIVE BOVINE ARRESTIN 1
Authors : Sander, C.L.; Palczewski, K.; Kiser, P.D.
Deposited on : 2020-08-14
Resolution : 2.50 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

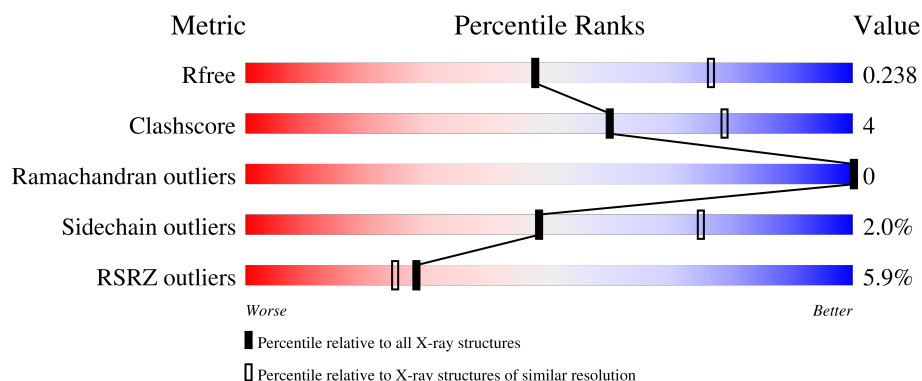
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	404	3% 82% 8% 9%
1	B	404	9% 81% 13% 5%
1	C	404	4% 80% 9% 10%
1	D	404	6% 83% 12% 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11919 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 S-arrestin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	S	0	1	0
			2913	1863	490	551	9			
1	B	383	Total	C	N	O	S	0	0	0
			2998	1912	499	577	10			
1	C	362	Total	C	N	O	S	0	0	0
			2847	1824	480	535	8			
1	D	383	Total	C	N	O	S	0	0	0
			3001	1911	500	580	10			

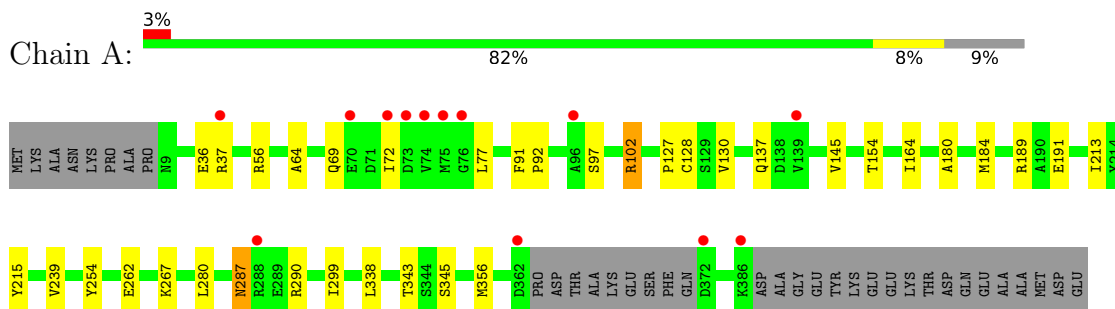
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	50	Total	O	0	0
			50	50		
2	B	27	Total	O	0	0
			27	27		
2	C	33	Total	O	0	0
			33	33		
2	D	50	Total	O	0	0
			50	50		

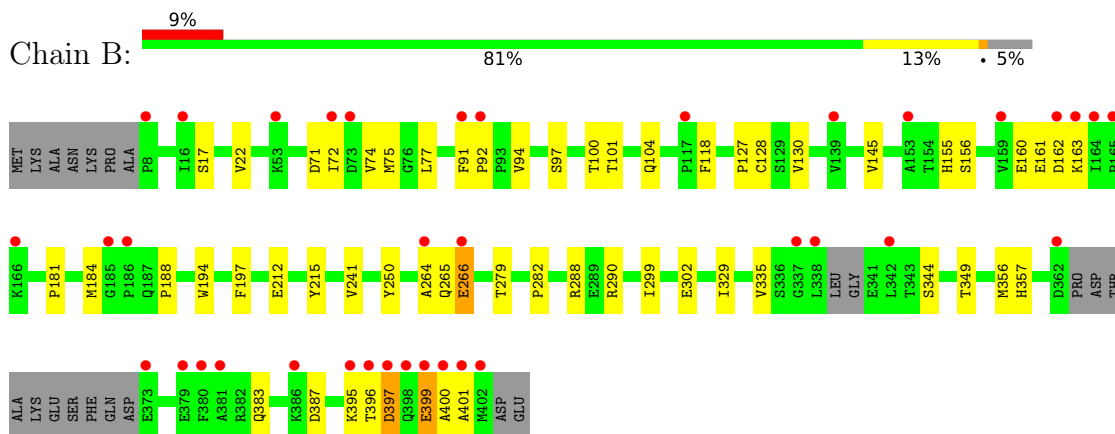
3 Residue-property plots [i](#)

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.

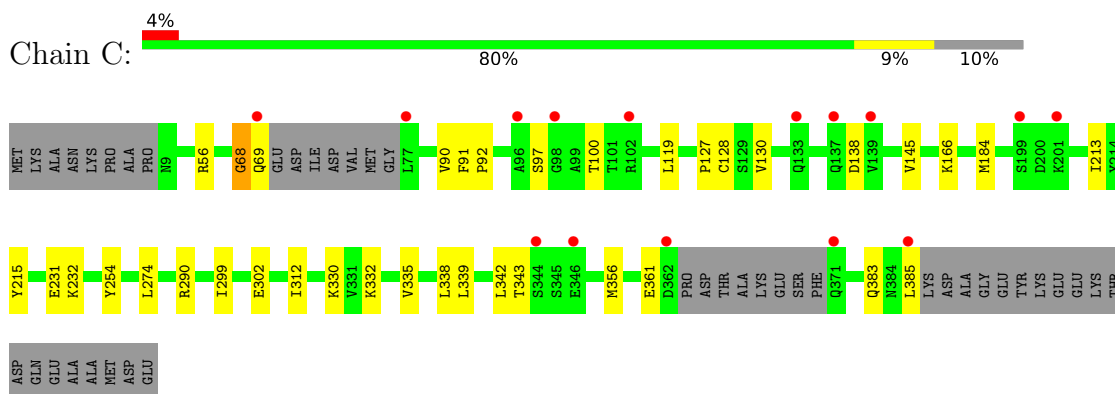
• Molecule 1: S-arrestin



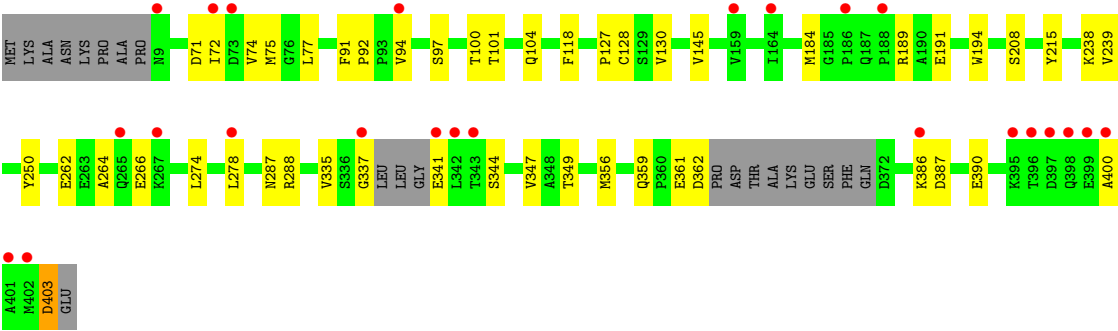
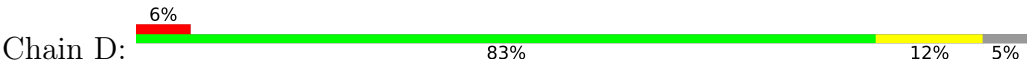
• Molecule 1: S-arrestin



• Molecule 1: S-arrestin



● Molecule 1: S-arrestin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	167.18Å 186.97Å 90.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.04 – 2.50 49.04 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.04-2.50) 99.9 (49.04-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.208 , 0.233 0.217 , 0.238	Depositor DCC
R_{free} test set	4928 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	68.8	Xtriage
Anisotropy	0.191	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.0	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	0.95	EDS
Total number of atoms	11919	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.64% 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 [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	1.01	0/2970	1.27	1/4027 (0.0%)
1	B	1.03	0/3056	1.29	1/4144 (0.0%)
1	C	0.98	0/2903	1.27	2/3936 (0.1%)
1	D	1.00	0/3058	1.29	1/4147 (0.0%)
All	All	1.00	0/11987	1.28	5/16254 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	90	VAL	CB-CA-C	6.84	117.08	111.06
1	C	68	GLY	CA-C-O	-5.49	118.65	122.22
1	A	287	ASN	CA-CB-CG	5.28	117.88	112.60
1	B	279	THR	CA-CB-OG1	-5.25	101.73	109.60
1	D	403	ASP	CA-CB-CG	5.10	117.70	112.60

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	2913	0	2959	24	0
1	B	2998	0	3008	40	0
1	C	2847	0	2901	17	0
1	D	3001	0	3002	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	50	0	0	0	0
2	B	27	0	0	1	0
2	C	33	0	0	0	0
2	D	50	0	0	1	0
All	All	11919	0	11870	101	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 4.

All (101) 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:B:401:ALA:HB1	1:C:138:ASP:OD1	1.78	0.83
1:B:397:ASP:OD1	1:B:397:ASP:N	2.11	0.83
1:A:254:TYR:CD2	1:D:400:ALA:HB2	2.20	0.75
1:B:399:GLU:OE2	1:B:399:GLU:HA	1.93	0.69
1:A:239:VAL:HG12	1:A:280:LEU:HD11	1.76	0.67
1:B:400:ALA:HB2	1:C:254:TYR:CD2	2.34	0.63
1:B:264:ALA:HB1	1:B:266:GLU:HG3	1.82	0.62
1:B:290:ARG:NH1	1:B:299:ILE:O	2.34	0.61
1:D:77:LEU:HD21	1:D:250:TYR:OH	2.00	0.61
1:D:287:ASN:HB2	2:D:517:HOH:O	2.03	0.58
1:B:94:VAL:HG12	1:B:94:VAL:O	2.05	0.57
1:D:184:MET:HE2	1:D:359:GLN:HE21	1.70	0.57
1:A:290:ARG:NH1	1:A:299:ILE:O	2.38	0.57
1:C:290:ARG:NH1	1:C:299:ILE:O	2.39	0.56
1:A:254:TYR:CD2	1:D:400:ALA:CB	2.89	0.56
1:A:215:TYR:HA	1:A:356:MET:O	2.06	0.56
1:C:215:TYR:HA	1:C:356:MET:O	2.06	0.56
1:B:215:TYR:HA	1:B:356:MET:O	2.06	0.56
1:A:137:GLN:HG3	1:D:403:ASP:CB	2.37	0.55
1:D:94:VAL:O	1:D:94:VAL:HG12	2.06	0.55
1:D:215:TYR:HA	1:D:356:MET:O	2.07	0.54
1:A:69:GLN:HB2	1:A:77:LEU:HD13	1.89	0.54
1:C:128:CYS:O	1:C:130:VAL:HG23	2.09	0.53
1:D:128:CYS:O	1:D:130:VAL:HG23	2.10	0.52
1:A:128:CYS:O	1:A:130:VAL:HG23	2.10	0.52
1:B:128:CYS:O	1:B:130:VAL:HG23	2.10	0.51
1:B:188:PRO:HG2	1:B:212:GLU:CD	2.35	0.51
1:D:390:GLU:HA	1:D:390:GLU:OE2	2.11	0.51
1:A:137:GLN:HG3	1:D:403:ASP:HB3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:GLU:OE2	1:B:163:LYS:HE3	2.10	0.51
1:B:75:MET:CE	1:B:77:LEU:HD12	2.42	0.50
1:B:194:TRP:CZ2	1:B:349:THR:HB	2.46	0.50
1:B:357:HIS:HE1	2:B:514:HOH:O	1.94	0.50
1:B:77:LEU:HD21	1:B:250:TYR:OH	2.12	0.49
1:D:239:VAL:HG21	1:D:278:LEU:HD13	1.94	0.49
1:C:184:MET:HB3	1:C:213:ILE:HD12	1.94	0.49
1:B:17:SER:HB3	1:B:22:VAL:CG1	2.43	0.48
1:A:36:GLU:C	1:A:37[B]:ARG:HG3	2.38	0.48
1:B:127:PRO:HG2	1:B:145:VAL:HG11	1.95	0.48
1:A:239:VAL:HG12	1:A:280:LEU:CD1	2.42	0.48
1:A:262:GLU:HB2	1:A:280:LEU:HD21	1.96	0.47
1:D:71:ASP:HB2	1:D:74:VAL:HB	1.96	0.47
1:A:184:MET:HB3	1:A:213:ILE:HD12	1.96	0.47
1:A:267:LYS:HD2	1:B:197:PHE:HB2	1.95	0.47
1:B:17:SER:HB3	1:B:22:VAL:HG12	1.97	0.47
1:C:338:LEU:HG	1:C:343:THR:HG22	1.95	0.47
1:D:337:GLY:O	1:D:341:GLU:N	2.48	0.46
1:A:64:ALA:HB3	1:A:72:ILE:HG21	1.96	0.46
1:D:127:PRO:HG2	1:D:145:VAL:HG11	1.96	0.46
1:D:194:TRP:CZ2	1:D:349:THR:HB	2.50	0.46
1:B:264:ALA:HB1	1:B:266:GLU:CG	2.46	0.46
1:A:191:GLU:O	1:D:77:LEU:HA	2.16	0.46
1:A:137:GLN:H	1:A:137:GLN:CD	2.23	0.45
1:A:189:ARG:HE	1:D:75:MET:HE2	1.80	0.45
1:D:264:ALA:HB1	1:D:266:GLU:OE2	2.17	0.45
1:B:181:PRO:HG2	1:B:184:MET:HG2	1.98	0.45
1:B:395:LYS:HE3	1:B:395:LYS:HB2	1.79	0.45
1:D:100:THR:HG22	1:D:101:THR:O	2.17	0.45
1:B:241:VAL:HG11	1:B:282:PRO:HB3	1.99	0.45
1:A:127:PRO:HG2	1:A:145:VAL:HG11	1.99	0.45
1:B:264:ALA:HB1	1:B:266:GLU:CD	2.43	0.45
1:D:288:ARG:NH1	1:D:387:ASP:OD1	2.50	0.44
1:B:335:VAL:O	1:B:344:SER:HB2	2.17	0.44
1:D:194:TRP:CD2	1:D:349:THR:HG21	2.52	0.44
1:B:71:ASP:HB2	1:B:74:VAL:HB	1.99	0.44
1:B:77:LEU:C	1:B:77:LEU:HD23	2.43	0.44
1:B:396:THR:O	1:B:397:ASP:C	2.60	0.44
1:C:91:PHE:HA	1:C:92:PRO:C	2.42	0.44
1:A:154:THR:HG22	1:A:164:ILE:HG12	2.00	0.43
1:B:396:THR:C	1:B:397:ASP:O	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:PHE:HA	1:A:92:PRO:C	2.43	0.43
1:B:288:ARG:NH1	1:B:387:ASP:OD1	2.51	0.43
1:D:335:VAL:O	1:D:344:SER:HB2	2.19	0.43
1:B:97:SER:OG	1:B:118:PHE:HA	2.19	0.43
1:B:91:PHE:HA	1:B:92:PRO:C	2.44	0.43
1:C:68:GLY:O	1:C:69:GLN:O	2.37	0.43
1:D:189:ARG:NH2	1:D:208:SER:OG	2.41	0.43
1:A:338:LEU:HG	1:A:343:THR:HG22	2.01	0.43
1:B:100:THR:HG22	1:B:104:GLN:HB2	2.01	0.42
1:C:100:THR:HG21	1:C:119:LEU:HD22	2.01	0.42
1:D:77:LEU:C	1:D:77:LEU:HD23	2.44	0.42
1:C:302:GLU:OE2	1:C:383:GLN:OE1	2.38	0.42
1:C:330:LYS:HD3	1:C:332:LYS:HE3	2.01	0.42
1:D:77:LEU:HD23	1:D:77:LEU:O	2.19	0.42
1:D:262:GLU:HB3	1:D:278:LEU:HD23	2.01	0.42
1:B:100:THR:HG22	1:B:101:THR:O	2.19	0.42
1:B:302:GLU:OE1	1:B:383:GLN:NE2	2.51	0.42
1:D:91:PHE:HA	1:D:92:PRO:C	2.44	0.42
1:B:265:GLN:HA	1:B:265:GLN:OE1	2.20	0.42
1:C:274:LEU:HD12	1:D:347:VAL:HG12	2.02	0.42
1:C:335:VAL:HB	1:C:338:LEU:HD22	2.00	0.42
1:A:180:ALA:HA	1:A:356:MET:SD	2.60	0.41
1:C:127:PRO:HG2	1:C:145:VAL:HG11	2.00	0.41
1:D:97:SER:OG	1:D:118:PHE:HA	2.19	0.41
1:C:232:LYS:HE3	1:C:342:LEU:HD12	2.03	0.41
1:D:100:THR:HG22	1:D:104:GLN:HB2	2.02	0.41
1:B:266:GLU:HG3	1:B:266:GLU:H	1.69	0.41
1:B:162:ASP:HA	1:C:339:LEU:HD22	2.01	0.41
1:B:194:TRP:CD2	1:B:349:THR:HG21	2.56	0.41
1:A:102:ARG:HD2	1:A:102:ARG:HA	1.90	0.40
1:B:241:VAL:HG22	1:B:329:ILE:HD13	2.03	0.40

There are no symmetry-related clashes.

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	366/404 (91%)	357 (98%)	9 (2%)	0	100	100
1	B	377/404 (93%)	366 (97%)	11 (3%)	0	100	100
1	C	356/404 (88%)	348 (98%)	8 (2%)	0	100	100
1	D	377/404 (93%)	366 (97%)	11 (3%)	0	100	100
All	All	1476/1616 (91%)	1437 (97%)	39 (3%)	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	328/359 (91%)	323 (98%)	5 (2%)	57	80
1	B	335/359 (93%)	328 (98%)	7 (2%)	47	74
1	C	321/359 (89%)	314 (98%)	7 (2%)	45	73
1	D	335/359 (93%)	328 (98%)	7 (2%)	47	74
All	All	1319/1436 (92%)	1293 (98%)	26 (2%)	48	75

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	A	97	SER
1	A	102	ARG
1	A	287	ASN
1	A	345	SER
1	B	72	ILE
1	B	155	HIS
1	B	156	SER
1	B	160	GLU
1	B	266	GLU

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Mol	Chain	Res	Type
1	B	397	ASP
1	B	399	GLU
1	C	56	ARG
1	C	97	SER
1	C	166	LYS
1	C	231	GLU
1	C	312	ILE
1	C	361	GLU
1	C	385	LEU
1	D	72	ILE
1	D	191	GLU
1	D	238	LYS
1	D	274	LEU
1	D	361	GLU
1	D	362	ASP
1	D	386	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	ASN
1	A	246	ASN
1	B	114	ASN
1	B	359	GLN
1	C	114	ASN
1	C	246	ASN
1	C	265	GLN
1	D	87	GLN
1	D	114	ASN
1	D	195	GLN
1	D	359	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

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	369/404 (91%)	0.12	13 (3%) 47 42	35, 64, 100, 138	1 (0%)
1	B	383/404 (94%)	0.64	37 (9%) 13 11	51, 79, 132, 151	0
1	C	362/404 (89%)	0.23	15 (4%) 41 37	54, 69, 104, 135	0
1	D	383/404 (94%)	0.49	24 (6%) 26 23	51, 78, 123, 145	0
All	All	1497/1616 (92%)	0.37	89 (5%) 28 24	35, 71, 122, 151	1 (0%)

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	396	THR	6.1
1	D	72	ILE	5.9
1	A	75	MET	5.8
1	B	402	MET	5.6
1	C	139	VAL	5.4
1	D	342	LEU	5.0
1	C	137	GLN	4.9
1	B	72	ILE	4.8
1	D	9	ASN	4.8
1	A	37[A]	ARG	4.5
1	B	362	ASP	4.5
1	B	8	PRO	4.3
1	B	342	LEU	4.2
1	B	337	GLY	4.0
1	D	159	VAL	4.0
1	B	159	VAL	3.7
1	B	401	ALA	3.7
1	D	337	GLY	3.7
1	B	380	PHE	3.7
1	D	343	THR	3.7
1	B	399	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	362	ASP	3.7
1	A	72	ILE	3.7
1	B	73	ASP	3.4
1	D	188	PRO	3.4
1	A	139	VAL	3.3
1	C	69	GLN	3.2
1	A	386	LYS	3.2
1	C	362	ASP	3.2
1	B	264	ALA	3.2
1	C	385	LEU	3.2
1	D	402	MET	3.2
1	D	400	ALA	3.1
1	C	346	GLU	3.1
1	B	92	PRO	3.1
1	B	397	ASP	3.1
1	B	162	ASP	3.1
1	B	396	THR	3.0
1	B	338	LEU	3.0
1	A	73	ASP	3.0
1	D	186	PRO	3.0
1	D	267	LYS	3.0
1	A	76	GLY	2.9
1	D	395	LYS	2.9
1	D	397	ASP	2.9
1	B	139	VAL	2.9
1	B	381	ALA	2.9
1	C	77	LEU	2.9
1	D	399	GLU	2.9
1	B	386	LYS	2.8
1	B	398	GLN	2.8
1	D	401	ALA	2.8
1	B	163	LYS	2.8
1	D	73	ASP	2.8
1	D	398	GLN	2.7
1	B	395	LYS	2.7
1	D	341	GLU	2.7
1	B	400	ALA	2.7
1	A	288	ARG	2.7
1	B	266	GLU	2.6
1	B	379	GLU	2.6
1	C	199	SER	2.6
1	D	94	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	96	ALA	2.5
1	D	386	LYS	2.5
1	A	74	VAL	2.5
1	C	371	GLN	2.4
1	A	70	GLU	2.3
1	C	96	ALA	2.2
1	B	16	ILE	2.2
1	B	53	LYS	2.2
1	D	278	LEU	2.2
1	C	102	ARG	2.2
1	B	186	PRO	2.2
1	C	201	LYS	2.2
1	B	117	PRO	2.2
1	B	165	PRO	2.2
1	D	265	GLN	2.1
1	B	153	ALA	2.1
1	C	344	SER	2.1
1	C	98	GLY	2.1
1	B	164	ILE	2.1
1	B	166	LYS	2.1
1	B	91	PHE	2.1
1	B	373	GLU	2.1
1	D	164	ILE	2.1
1	A	372	ASP	2.1
1	C	133	GLN	2.0
1	B	185	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.