



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:46 AM UTC

PDB ID : 1SUJ / pdb_00001suj
Title : X-ray crystal structure of ambystoma tigrinum cone arrestin
Authors : Sutton, R.B.; Navarro, J.
Deposited on : 2004-03-26
Resolution : 2.38 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

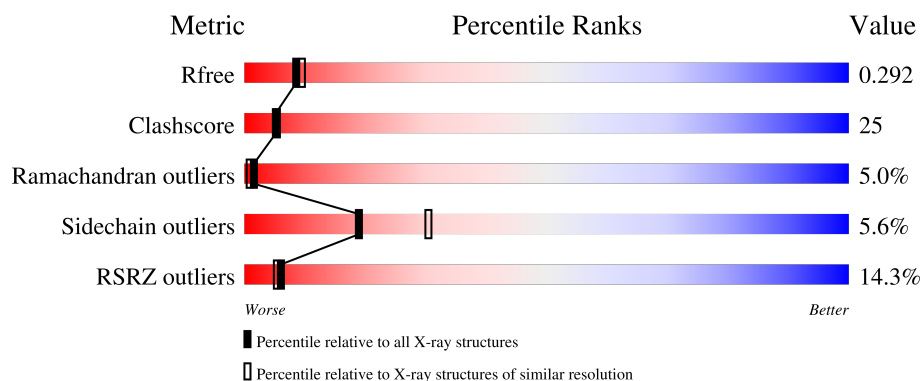
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	392	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2945 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 cone arrestin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2846	1818	481	533	14			

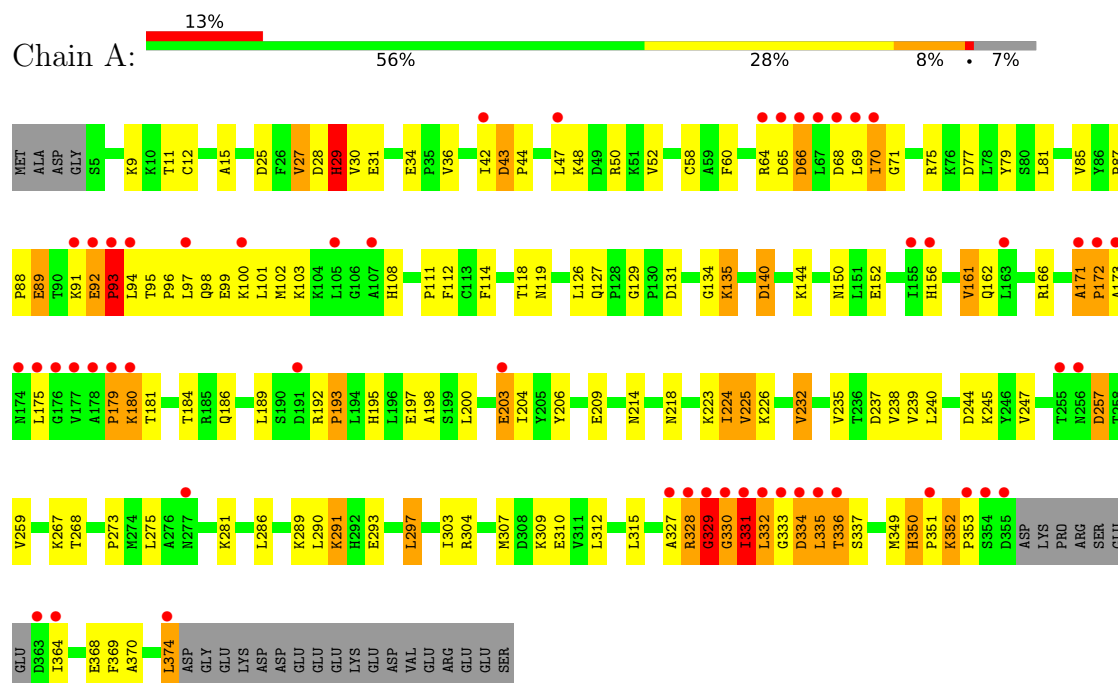
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	99	Total	O	0	0
			99	99		

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: cone arrestin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	70.76Å 75.05Å 79.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.22 – 2.38 27.22 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.9 (27.22-2.38) 99.8 (27.22-2.38)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.36Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.246 , 0.297 0.244 , 0.292	Depositor DCC
R_{free} test set	1640 reflections (9.33%)	wwPDB-VP
Wilson B-factor (Å ²)	37.0	Xtriage
Anisotropy	0.251	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.022 for -h,l,k	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2945	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.85% 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	0.41	0/2897	0.99	21/3920 (0.5%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	ALA	N-CA-C	-8.02	103.39	113.01
1	A	29	HIS	N-CA-C	-7.88	94.01	110.80
1	A	127	GLN	CA-C-N	6.81	126.84	119.89
1	A	127	GLN	C-N-CA	6.81	126.84	119.89
1	A	350	HIS	CA-C-N	6.43	127.88	119.84
1	A	350	HIS	C-N-CA	6.43	127.88	119.84
1	A	129	GLY	CA-C-N	5.97	126.39	119.47
1	A	129	GLY	C-N-CA	5.97	126.39	119.47
1	A	336	THR	N-CA-C	-5.90	102.48	110.68
1	A	329	GLY	N-CA-C	-5.80	99.42	113.18
1	A	238	VAL	N-CA-C	-5.75	99.42	108.23
1	A	171	ALA	CA-C-N	5.74	127.01	119.84
1	A	171	ALA	C-N-CA	5.74	127.01	119.84
1	A	27	VAL	N-CA-C	5.73	116.83	108.58
1	A	92	GLU	N-CA-C	5.73	119.44	109.82
1	A	232	VAL	N-CA-C	-5.58	99.70	107.80
1	A	43	ASP	CA-C-N	5.57	125.03	119.24
1	A	43	ASP	C-N-CA	5.57	125.03	119.24
1	A	224	ILE	N-CA-C	5.40	115.67	108.11
1	A	337	SER	N-CA-C	5.26	117.32	110.53
1	A	193	PRO	N-CA-C	5.05	119.12	111.19

There are no chirality outliers.

There are no planarity outliers.

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	2846	0	2940	142	0
2	A	99	0	0	3	0
All	All	2945	0	2940	142	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 25.

All (142) 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:172:PRO:HB2	1:A:175:LEU:HG	1.41	1.03
1:A:135:LYS:HZ2	1:A:135:LYS:HB3	1.26	0.98
1:A:331:ILE:HB	1:A:332:LEU:HD23	1.49	0.94
1:A:327:ALA:HB3	1:A:330:GLY:HA2	1.49	0.93
1:A:291:LYS:HE2	1:A:291:LYS:H	1.32	0.93
1:A:100:LYS:HB3	1:A:364:ILE:HD11	1.57	0.84
1:A:225:VAL:HG13	1:A:259:VAL:HB	1.57	0.84
1:A:44:PRO:HG2	1:A:108:HIS:CE1	2.13	0.84
1:A:291:LYS:H	1:A:291:LYS:CE	1.91	0.83
1:A:289:LYS:HB3	1:A:291:LYS:HE3	1.61	0.83
1:A:303:ILE:HD13	1:A:309:LYS:HD2	1.63	0.81
1:A:329:GLY:O	1:A:331:ILE:HG13	1.81	0.80
1:A:327:ALA:C	1:A:329:GLY:H	1.91	0.79
1:A:144:LYS:HG2	1:A:162:GLN:HG2	1.67	0.76
1:A:235:VAL:HG13	1:A:245:LYS:HE3	1.66	0.75
1:A:224:ILE:HB	1:A:330:GLY:HA3	1.68	0.75
1:A:286:LEU:HD11	1:A:290:LEU:HD23	1.68	0.74
1:A:69:LEU:HG	1:A:70:ILE:HG22	1.68	0.74
1:A:304:ARG:O	1:A:307:MET:HG2	1.87	0.74
1:A:232:VAL:HG21	1:A:273:PRO:HG3	1.71	0.72
1:A:144:LYS:HG2	1:A:162:GLN:CG	2.20	0.72
1:A:327:ALA:O	1:A:329:GLY:N	2.25	0.70
1:A:226:LYS:HE3	1:A:330:GLY:HA2	1.74	0.69
1:A:65:ASP:OD2	1:A:69:LEU:HD22	1.92	0.69
1:A:99:GLU:O	1:A:103:LYS:HG2	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:PRO:O	1:A:352:LYS:HB2	1.92	0.68
1:A:203:GLU:HB3	1:A:204:ILE:HD12	1.77	0.67
1:A:327:ALA:HB3	1:A:330:GLY:CA	2.23	0.66
1:A:27:VAL:HG21	1:A:353:PRO:HG3	1.75	0.66
1:A:226:LYS:HE3	1:A:330:GLY:CA	2.26	0.66
1:A:275:LEU:HB2	1:A:297:LEU:HD13	1.77	0.66
1:A:28:ASP:OD1	1:A:29:HIS:O	2.14	0.66
1:A:333:GLY:C	1:A:335:LEU:H	2.04	0.65
1:A:291:LYS:HE2	1:A:291:LYS:N	2.09	0.64
1:A:328:ARG:HG3	1:A:336:THR:O	1.97	0.64
1:A:281:LYS:CB	1:A:290:LEU:HD22	2.29	0.63
1:A:27:VAL:HG21	1:A:353:PRO:CG	2.28	0.62
1:A:330:GLY:C	1:A:331:ILE:HG13	2.24	0.62
1:A:12:CYS:HB3	1:A:161:VAL:HG22	1.82	0.62
1:A:286:LEU:CD1	1:A:290:LEU:HD23	2.30	0.61
1:A:171:ALA:HB2	1:A:349:MET:HG3	1.83	0.61
1:A:119:ASN:OD1	1:A:304:ARG:HD2	2.01	0.61
1:A:172:PRO:HD2	1:A:175:LEU:HD11	1.83	0.61
1:A:334:ASP:C	1:A:336:THR:H	2.09	0.61
1:A:42:ILE:HG22	1:A:43:ASP:N	2.15	0.60
1:A:327:ALA:C	1:A:329:GLY:N	2.60	0.60
1:A:329:GLY:O	1:A:331:ILE:N	2.35	0.60
1:A:237:ASP:OD2	1:A:245:LYS:HD3	2.01	0.60
1:A:281:LYS:HB3	1:A:290:LEU:CD2	2.32	0.59
1:A:100:LYS:CB	1:A:364:ILE:HD11	2.30	0.58
1:A:267:LYS:HG2	1:A:268:THR:N	2.18	0.58
1:A:286:LEU:HD11	1:A:290:LEU:CD2	2.32	0.58
1:A:206:TYR:O	1:A:209:GLU:HB3	2.04	0.57
1:A:48:LYS:O	1:A:48:LYS:HG3	2.04	0.57
1:A:245:LYS:HE2	1:A:247:VAL:CG2	2.34	0.57
1:A:206:TYR:HA	1:A:349:MET:O	2.05	0.57
1:A:135:LYS:HB3	1:A:135:LYS:NZ	2.09	0.57
1:A:351:PRO:O	1:A:352:LYS:CB	2.53	0.56
1:A:144:LYS:HE2	1:A:162:GLN:CD	2.30	0.56
1:A:281:LYS:HB3	1:A:290:LEU:HD22	1.86	0.56
1:A:327:ALA:CB	1:A:330:GLY:HA2	2.27	0.56
1:A:331:ILE:HB	1:A:332:LEU:CD2	2.30	0.56
1:A:42:ILE:HG21	1:A:47:LEU:HD22	1.88	0.56
1:A:30:VAL:HB	1:A:172:PRO:HG3	1.86	0.56
1:A:331:ILE:HG22	1:A:332:LEU:H	1.70	0.56
1:A:75:ARG:HG2	1:A:75:ARG:HH11	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:VAL:CG1	1:A:259:VAL:HB	2.31	0.55
1:A:31:GLU:HB2	2:A:486:HOH:O	2.07	0.54
1:A:293:GLU:HB2	1:A:374:LEU:HB2	1.89	0.53
1:A:118:THR:O	1:A:119:ASN:HB2	2.09	0.53
1:A:27:VAL:O	1:A:34:GLU:HG3	2.09	0.53
1:A:42:ILE:HD13	1:A:52:VAL:HG21	1.90	0.53
1:A:119:ASN:ND2	1:A:304:ARG:NH1	2.57	0.52
1:A:245:LYS:HE2	1:A:247:VAL:HG22	1.91	0.52
1:A:235:VAL:CG1	1:A:245:LYS:HE3	2.36	0.52
1:A:28:ASP:C	1:A:29:HIS:O	2.51	0.51
1:A:25:ASP:CG	1:A:166:ARG:HH21	2.19	0.51
1:A:184:THR:HB	1:A:197:GLU:HG3	1.92	0.51
1:A:95:THR:OG1	1:A:98:GLN:HG3	2.12	0.50
1:A:333:GLY:C	1:A:335:LEU:N	2.70	0.50
1:A:64:ARG:NH2	1:A:68:ASP:OD1	2.45	0.50
1:A:112:PHE:CD1	1:A:112:PHE:C	2.89	0.49
1:A:303:ILE:HD11	1:A:315:LEU:HD22	1.93	0.49
1:A:368:GLU:HG3	2:A:406:HOH:O	2.12	0.49
1:A:328:ARG:HD2	1:A:336:THR:HB	1.93	0.49
1:A:181:THR:HG21	2:A:447:HOH:O	2.11	0.49
1:A:281:LYS:HB2	1:A:290:LEU:HD22	1.94	0.48
1:A:65:ASP:O	1:A:66:ASP:C	2.57	0.48
1:A:179:PRO:O	1:A:180:LYS:HG3	2.13	0.47
1:A:186:GLN:HG2	1:A:195:HIS:ND1	2.29	0.47
1:A:30:VAL:CG2	1:A:172:PRO:HG3	2.44	0.47
1:A:198:ALA:HA	1:A:214:ASN:O	2.15	0.47
1:A:126:LEU:C	1:A:126:LEU:HD23	2.39	0.47
1:A:144:LYS:HG2	1:A:162:GLN:HG3	1.94	0.47
1:A:150:ASN:OD1	1:A:152:GLU:HB2	2.15	0.47
1:A:240:LEU:HD23	1:A:240:LEU:HA	1.78	0.47
1:A:333:GLY:O	1:A:335:LEU:N	2.41	0.47
1:A:330:GLY:C	1:A:331:ILE:CG1	2.88	0.47
1:A:94:LEU:HD21	1:A:102:MET:SD	2.56	0.46
1:A:96:PRO:O	1:A:99:GLU:HB2	2.16	0.45
1:A:331:ILE:C	1:A:332:LEU:HD23	2.42	0.45
1:A:257:ASP:OD1	1:A:267:LYS:NZ	2.45	0.45
1:A:92:GLU:HB3	1:A:93:PRO:HD2	1.97	0.45
1:A:42:ILE:CG2	1:A:43:ASP:N	2.80	0.44
1:A:75:ARG:HG2	1:A:75:ARG:NH1	2.32	0.44
1:A:369:PHE:O	1:A:370:ALA:C	2.61	0.44
1:A:75:ARG:NH1	1:A:77:ASP:OD1	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ILE:HG22	1:A:43:ASP:H	1.83	0.44
1:A:102:MET:CE	1:A:111:PRO:HD3	2.48	0.44
1:A:36:VAL:HB	1:A:114:PHE:HB2	2.00	0.44
1:A:134:GLY:HA3	1:A:244:ASP:OD2	2.18	0.43
1:A:239:VAL:O	1:A:310:GLU:HG2	2.18	0.43
1:A:97:LEU:C	1:A:97:LEU:HD12	2.43	0.43
1:A:275:LEU:HD11	1:A:281:LYS:HD3	1.99	0.43
1:A:352:LYS:HG2	1:A:353:PRO:CD	2.49	0.43
1:A:11:THR:CG2	1:A:15:ALA:HA	2.48	0.43
1:A:64:ARG:HH11	1:A:64:ARG:HG2	1.83	0.43
1:A:352:LYS:HG2	1:A:353:PRO:HD2	2.01	0.43
1:A:331:ILE:O	1:A:332:LEU:C	2.62	0.43
1:A:52:VAL:O	1:A:85:VAL:HG22	2.18	0.42
1:A:189:LEU:HB3	1:A:223:LYS:HD2	2.00	0.42
1:A:30:VAL:CB	1:A:172:PRO:HG3	2.49	0.42
1:A:58:CYS:HA	1:A:140:ASP:O	2.20	0.42
1:A:47:LEU:CD1	1:A:50:ARG:HB2	2.49	0.42
1:A:203:GLU:C	1:A:204:ILE:HD12	2.45	0.42
1:A:94:LEU:HG	1:A:98:GLN:HB2	2.02	0.42
1:A:195:HIS:HB2	1:A:218:ASN:HB3	2.02	0.42
1:A:235:VAL:HG13	1:A:245:LYS:CE	2.43	0.42
1:A:97:LEU:O	1:A:101:LEU:HG	2.19	0.42
1:A:60:PHE:CD2	1:A:240:LEU:HD12	2.55	0.42
1:A:171:ALA:HA	1:A:172:PRO:HD3	1.94	0.42
1:A:334:ASP:C	1:A:336:THR:N	2.76	0.42
1:A:47:LEU:HD21	1:A:52:VAL:HG23	2.02	0.41
1:A:94:LEU:HD23	1:A:99:GLU:OE1	2.20	0.41
1:A:87:PRO:HA	1:A:88:PRO:HD3	1.87	0.41
1:A:225:VAL:O	1:A:225:VAL:HG22	2.20	0.41
1:A:303:ILE:HD13	1:A:309:LYS:CD	2.44	0.41
1:A:79:TYR:OH	1:A:81:LEU:HD11	2.20	0.41
1:A:58:CYS:HB3	1:A:79:TYR:HB3	2.02	0.40
1:A:11:THR:HG22	1:A:15:ALA:HA	2.04	0.40
1:A:192:ARG:HA	1:A:193:PRO:HD3	1.83	0.40
1:A:48:LYS:HE3	1:A:48:LYS:HB2	1.83	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	359/392 (92%)	320 (89%)	21 (6%)	18 (5%)	1 1

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	ASP
1	A	70	ILE
1	A	71	GLY
1	A	180	LYS
1	A	328	ARG
1	A	330	GLY
1	A	332	LEU
1	A	352	LYS
1	A	329	GLY
1	A	93	PRO
1	A	179	PRO
1	A	257	ASP
1	A	334	ASP
1	A	29	HIS
1	A	172	PRO
1	A	335	LEU
1	A	89	GLU
1	A	331	ILE

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	324/350 (93%)	306 (94%)	18 (6%)	19	30

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LYS
1	A	89	GLU
1	A	91	LYS
1	A	93	PRO
1	A	131	ASP
1	A	135	LYS
1	A	140	ASP
1	A	156	HIS
1	A	161	VAL
1	A	200	LEU
1	A	203	GLU
1	A	225	VAL
1	A	291	LYS
1	A	297	LEU
1	A	312	LEU
1	A	331	ILE
1	A	350	HIS
1	A	374	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	HIS
1	A	162	GLN
1	A	186	GLN
1	A	195	HIS
1	A	214	ASN
1	A	262	ASN
1	A	296	ASN
1	A	372	GLN

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](#)

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	363/392 (92%)	0.80	52 (14%) 6 5	16, 34, 79, 91	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	330	GLY	10.0
1	A	65	ASP	9.2
1	A	173	ALA	7.8
1	A	178	ALA	7.7
1	A	67	LEU	7.0
1	A	69	LEU	6.5
1	A	332	LEU	6.3
1	A	177	VAL	6.1
1	A	174	ASN	5.2
1	A	179	PRO	4.9
1	A	175	LEU	4.9
1	A	333	GLY	4.8
1	A	331	ILE	4.8
1	A	374	LEU	4.8
1	A	328	ARG	4.2
1	A	66	ASP	4.1
1	A	70	ILE	4.0
1	A	42	ILE	3.5
1	A	64	ARG	3.2
1	A	91	LYS	3.2
1	A	92	GLU	3.1
1	A	163	LEU	3.1
1	A	68	ASP	3.0
1	A	107	ALA	3.0
1	A	171	ALA	2.9
1	A	94	LEU	2.9
1	A	336	THR	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	176	GLY	2.8
1	A	334	ASP	2.7
1	A	363	ASP	2.7
1	A	180	LYS	2.6
1	A	355	ASP	2.6
1	A	203	GLU	2.5
1	A	329	GLY	2.5
1	A	277	ASN	2.4
1	A	353	PRO	2.4
1	A	255	THR	2.4
1	A	47	LEU	2.4
1	A	354	SER	2.3
1	A	327	ALA	2.3
1	A	256	ASN	2.3
1	A	364	ILE	2.2
1	A	105	LEU	2.2
1	A	93	PRO	2.2
1	A	351	PRO	2.2
1	A	156	HIS	2.1
1	A	335	LEU	2.1
1	A	191	ASP	2.1
1	A	100	LYS	2.0
1	A	97	LEU	2.0
1	A	155	ILE	2.0
1	A	172	PRO	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.