



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:58 AM UTC

PDB ID : 6IVP / pdb_00006ivp
Title : Crystal structure of a membrane protein P262A
Authors : Kittredge, A.; Fukuda, F.; Zhang, Y.; Yang, T.
Deposited on : 2018-12-04
Resolution : 3.80 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

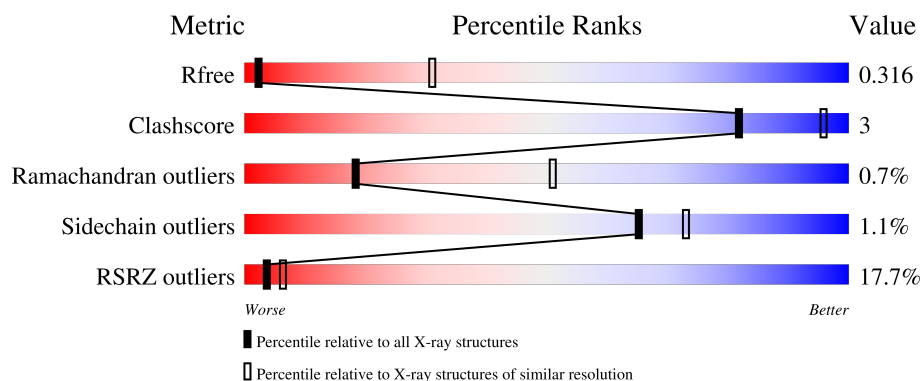
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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	1065 (3.96-3.64)
Clashscore	190562	1012 (3.94-3.66)
Ramachandran outliers	187476	1048 (3.96-3.64)
Sidechain outliers	187428	1043 (3.96-3.64)
RSRZ outliers	180081	1064 (3.96-3.64)

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	297	9% 86% 6% 8%
1	B	297	11% 81% 9% 9%
1	C	297	19% 82% 8% 9%
1	D	297	26% 83% 7% 10%
1	E	297	15% 82% 8% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10751 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 bestrophin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2175	1408	368	390	9			
1	B	269	Total	C	N	O	S	0	0	0
			2143	1387	364	383	9			
1	C	269	Total	C	N	O	S	0	0	0
			2138	1383	364	382	9			
1	D	268	Total	C	N	O	S	0	0	0
			2127	1375	361	382	9			
1	E	269	Total	C	N	O	S	0	0	0
			2143	1387	364	383	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP W9BH30
A	-1	ASN	-	expression tag	UNP W9BH30
A	0	ALA	-	expression tag	UNP W9BH30
A	262	ALA	PRO	engineered mutation	UNP W9BH30
B	-2	SER	-	expression tag	UNP W9BH30
B	-1	ASN	-	expression tag	UNP W9BH30
B	0	ALA	-	expression tag	UNP W9BH30
B	262	ALA	PRO	engineered mutation	UNP W9BH30
C	-2	SER	-	expression tag	UNP W9BH30
C	-1	ASN	-	expression tag	UNP W9BH30
C	0	ALA	-	expression tag	UNP W9BH30
C	262	ALA	PRO	engineered mutation	UNP W9BH30
D	-2	SER	-	expression tag	UNP W9BH30
D	-1	ASN	-	expression tag	UNP W9BH30
D	0	ALA	-	expression tag	UNP W9BH30
D	262	ALA	PRO	engineered mutation	UNP W9BH30
E	-2	SER	-	expression tag	UNP W9BH30
E	-1	ASN	-	expression tag	UNP W9BH30
E	0	ALA	-	expression tag	UNP W9BH30

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Chain	Residue	Modelled	Actual	Comment	Reference
E	262	ALA	PRO	engineered mutation	UNP W9BH30

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Zn 3 3	0	0
2	B	3	Total Zn 3 3	0	0
2	C	4	Total Zn 4 4	0	0
2	D	3	Total Zn 3 3	0	0
2	E	1	Total Zn 1 1	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Cl 2 2	0	0
3	B	2	Total Cl 2 2	0	0
3	C	2	Total Cl 2 2	0	0
3	D	3	Total Cl 3 3	0	0
3	E	1	Total Cl 1 1	0	0

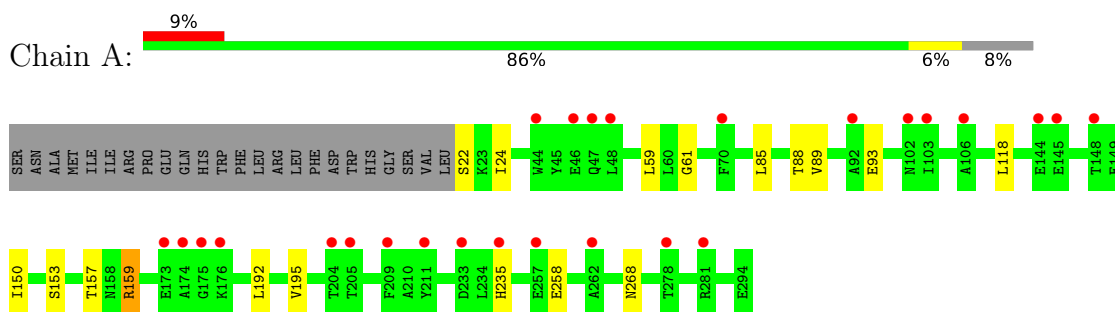
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total O 1 1	0	0

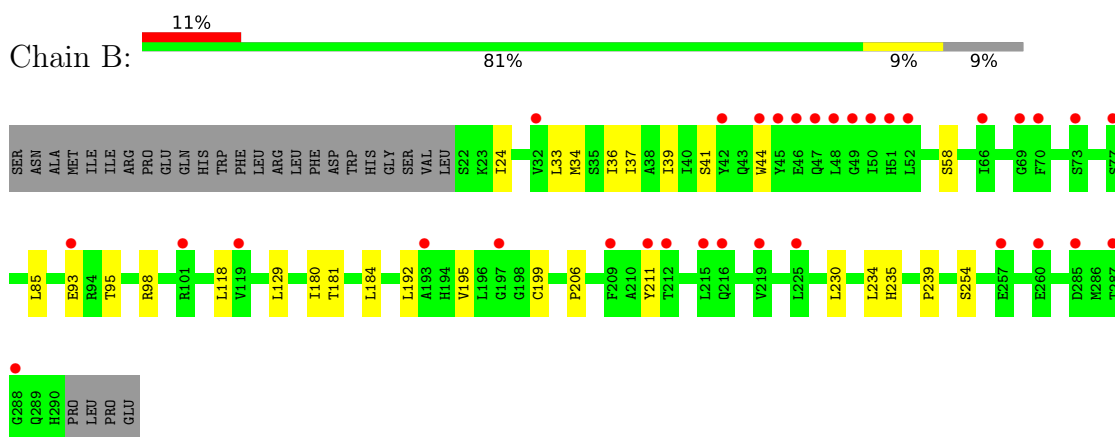
3 Residue-property plots [i](#)

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.

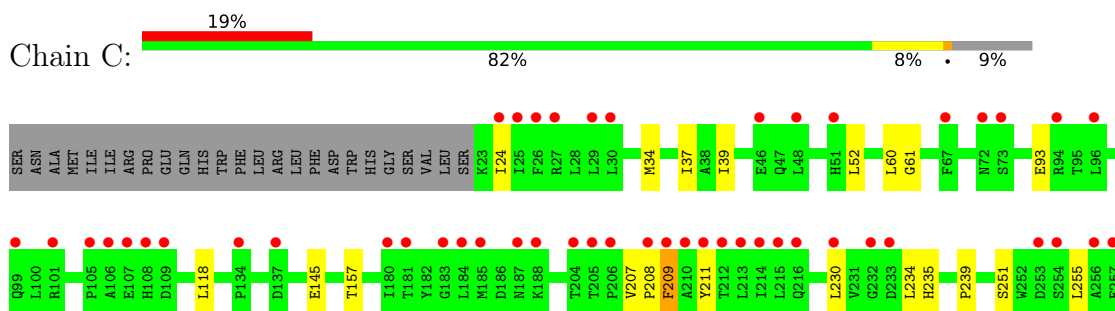
- Molecule 1: bestrophin

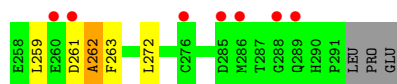


- Molecule 1: bestrophin

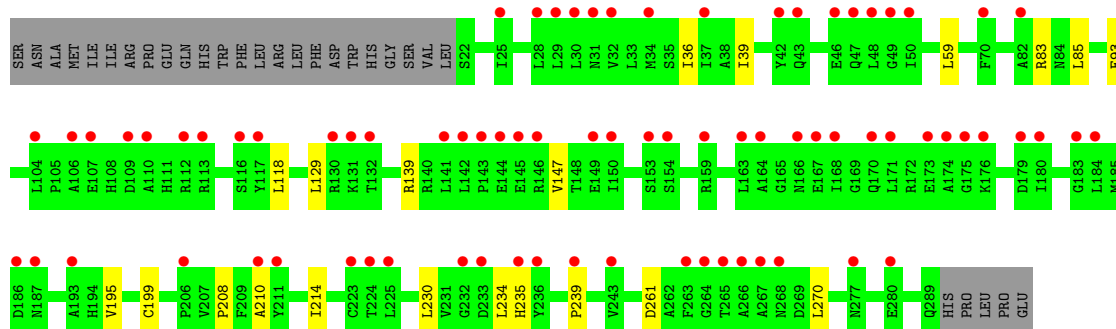
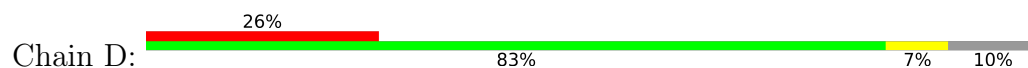


- Molecule 1: bestrophin

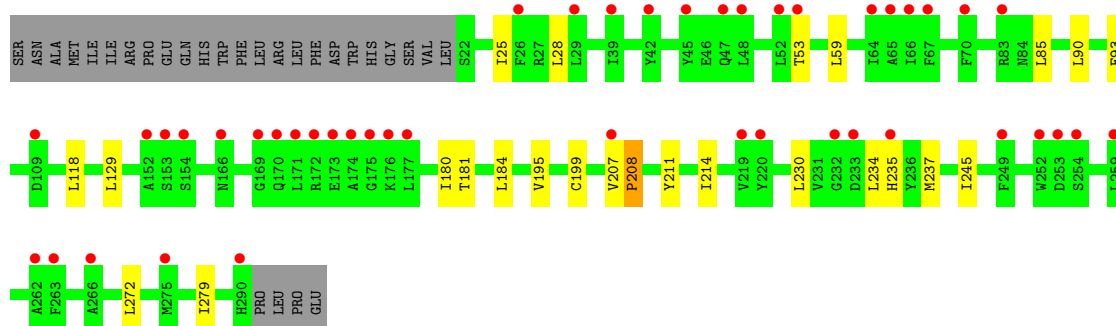
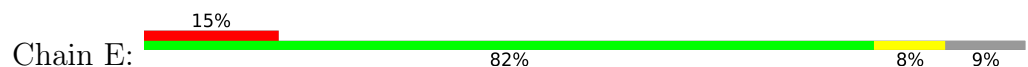




- Molecule 1: bestrophin



- Molecule 1: bestrophin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	114.48Å 162.41Å 161.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.01 – 3.80 50.01 – 3.80	Depositor EDS
% Data completeness (in resolution range)	97.5 (50.01-3.80) 97.5 (50.01-3.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 3.77Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.266 , 0.315 0.266 , 0.316	Depositor DCC
R_{free} test set	1526 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	149.8	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.22 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.026 for -h,l,k	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	10751	wwPDB-VP
Average B, all atoms (Å ²)	159.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.27% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, ZN

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	0.55	0/2223	1.02	0/3025
1	B	0.55	0/2189	1.02	0/2978
1	C	0.54	0/2184	1.06	2/2973 (0.1%)
1	D	0.54	0/2171	1.03	1/2954 (0.0%)
1	E	0.54	0/2189	1.02	0/2978
All	All	0.54	0/10956	1.03	3/14908 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	207	VAL	CA-C-N	5.80	127.09	119.84
1	C	207	VAL	C-N-CA	5.80	127.09	119.84
1	D	261	ASP	N-CA-C	5.51	116.98	110.97

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	2175	0	2217	10	0
1	B	2143	0	2186	14	0
1	C	2138	0	2181	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2127	0	2172	12	0
1	E	2143	0	2186	14	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	4	0	0	0	0
2	D	3	0	0	0	0
2	E	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	3	0	0	0	0
3	E	1	0	0	0	0
4	C	1	0	0	0	0
All	All	10751	0	10942	55	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 3.

All (55) 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:C:261:ASP:HA	1:C:262:ALA:HB2	1.37	1.03
1:C:261:ASP:HA	1:C:262:ALA:CB	2.16	0.67
1:E:207:VAL:HG23	1:E:208:PRO:HD3	1.79	0.65
1:E:90:LEU:HA	1:E:279:ILE:HD13	1.84	0.58
1:D:139:ARG:HG3	1:D:147:VAL:HG11	1.87	0.56
1:B:129:LEU:HD21	1:B:199:CYS:HB3	1.89	0.54
1:A:22:SER:HB3	1:A:24:ILE:HG22	1.91	0.53
1:E:129:LEU:HD21	1:E:199:CYS:HB3	1.90	0.52
1:D:230:LEU:HB3	1:D:234:LEU:HD12	1.90	0.52
1:C:39:ILE:HD11	1:C:239:PRO:HD3	1.93	0.51
1:E:230:LEU:HB3	1:E:234:LEU:HD12	1.93	0.51
1:E:211:TYR:HA	1:E:214:ILE:HB	1.93	0.50
1:A:89:VAL:HG22	1:A:195:VAL:HG11	1.93	0.50
1:C:230:LEU:HB3	1:C:234:LEU:HD12	1.93	0.50
1:D:93:GLU:HG2	1:D:118:LEU:HD22	1.93	0.50
1:A:153:SER:HB2	1:A:159:ARG:HD2	1.93	0.50
1:B:85:LEU:HB3	1:B:195:VAL:HG13	1.94	0.50
1:A:93:GLU:HG2	1:A:118:LEU:HD22	1.94	0.49
1:B:39:ILE:HD11	1:B:239:PRO:HD3	1.95	0.49
1:D:85:LEU:HD22	1:D:195:VAL:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:LEU:HB3	1:B:234:LEU:HD12	1.94	0.48
1:B:34:MET:HA	1:B:37:ILE:HD12	1.96	0.48
1:C:61:GLY:HA3	1:D:59:LEU:HD11	1.96	0.48
1:D:83:ARG:HG2	1:D:270:LEU:HD21	1.95	0.48
1:B:85:LEU:HD22	1:B:195:VAL:HA	1.96	0.47
1:C:251:SER:HB3	1:D:214:ILE:HG12	1.96	0.47
1:B:93:GLU:HG2	1:B:118:LEU:HD22	1.96	0.47
1:C:93:GLU:HG2	1:C:118:LEU:HD22	1.97	0.47
1:E:93:GLU:HG2	1:E:118:LEU:HD22	1.96	0.47
1:B:95:THR:HG22	1:B:98:ARG:HE	1.79	0.46
1:D:39:ILE:HD11	1:D:239:PRO:HD3	1.99	0.45
1:D:36:ILE:HA	1:D:39:ILE:HD12	1.98	0.45
1:C:34:MET:HA	1:C:37:ILE:HD12	1.98	0.45
1:C:263:PHE:HB3	1:C:272:LEU:HD11	1.99	0.45
1:C:60:LEU:HB2	1:E:245:ILE:HG12	1.99	0.44
1:B:33:LEU:HA	1:B:36:ILE:HD12	1.99	0.44
1:C:60:LEU:HD13	1:E:245:ILE:HA	2.00	0.44
1:E:85:LEU:HD22	1:E:195:VAL:HA	1.99	0.44
1:C:259:LEU:HD21	1:D:208:PRO:HG2	2.01	0.43
1:A:258:GLU:HB3	1:A:268:ASN:HB2	2.00	0.43
1:E:25:ILE:HA	1:E:28:LEU:HD12	2.00	0.43
1:B:24:ILE:HD11	1:B:254:SER:HB3	2.01	0.42
1:A:150:ILE:HA	1:A:159:ARG:HG2	2.02	0.42
1:B:181:THR:HA	1:B:184:LEU:HD12	2.01	0.42
1:B:93:GLU:HG3	1:B:192:LEU:HD21	2.00	0.42
1:E:181:THR:HA	1:E:184:LEU:HD12	2.02	0.42
1:A:93:GLU:HG3	1:A:192:LEU:HD21	2.02	0.42
1:A:61:GLY:HA3	1:E:59:LEU:HD11	2.01	0.42
1:A:85:LEU:HD22	1:A:195:VAL:HA	2.02	0.41
1:E:85:LEU:HB3	1:E:195:VAL:HG13	2.03	0.41
1:D:129:LEU:HD21	1:D:199:CYS:HB3	2.01	0.41
1:A:59:LEU:HD13	1:B:58:SER:HA	2.03	0.40
1:B:41:SER:HA	1:B:44:TRP:HD1	1.87	0.40
1:C:255:LEU:HD13	1:D:210:ALA:HB3	2.03	0.40
1:C:52:LEU:HD11	1:E:237:MET:HE2	2.04	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	271/297 (91%)	266 (98%)	4 (2%)	1 (0%)	30	62
1	B	267/297 (90%)	262 (98%)	3 (1%)	2 (1%)	18	51
1	C	267/297 (90%)	259 (97%)	4 (2%)	4 (2%)	8	36
1	D	266/297 (90%)	263 (99%)	2 (1%)	1 (0%)	30	62
1	E	267/297 (90%)	259 (97%)	6 (2%)	2 (1%)	18	51
All	All	1338/1485 (90%)	1309 (98%)	19 (1%)	10 (1%)	18	51

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	209	PHE
1	C	262	ALA
1	A	235	HIS
1	C	235	HIS
1	B	235	HIS
1	D	235	HIS
1	E	235	HIS
1	B	206	PRO
1	C	208	PRO
1	E	208	PRO

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	237/259 (92%)	234 (99%)	3 (1%)	61	71
1	B	233/259 (90%)	231 (99%)	2 (1%)	70	74
1	C	232/259 (90%)	227 (98%)	5 (2%)	45	63
1	D	231/259 (89%)	231 (100%)	0	100	100
1	E	233/259 (90%)	230 (99%)	3 (1%)	61	71
All	All	1166/1295 (90%)	1153 (99%)	13 (1%)	65	73

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	THR
1	A	157	THR
1	A	159	ARG
1	B	180	ILE
1	B	211	TYR
1	C	24	ILE
1	C	145	GLU
1	C	157	THR
1	C	209	PHE
1	C	211	TYR
1	E	53	THR
1	E	180	ILE
1	E	272	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	72	ASN
1	A	99	GLN
1	A	158	ASN
1	B	99	GLN
1	B	102	ASN
1	B	273	ASN
1	B	290	HIS
1	C	31	ASN
1	C	99	GLN
1	C	102	ASN
1	C	128	GLN
1	C	166	ASN

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Mol	Chain	Res	Type
1	C	273	ASN
1	D	31	ASN
1	D	47	GLN
1	D	51	HIS
1	D	102	ASN
1	D	166	ASN
1	D	187	ASN
1	D	268	ASN
1	D	277	ASN
1	D	282	ASN
1	D	289	GLN
1	E	31	ASN
1	E	47	GLN
1	E	99	GLN
1	E	102	ASN
1	E	166	ASN
1	E	268	ASN

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](#)

Of 24 ligands modelled in this entry, 24 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	273/297 (91%)	0.71	26 (9%) 14 15	104, 149, 206, 261	0
1	B	269/297 (90%)	0.90	33 (12%) 8 12	108, 163, 210, 227	0
1	C	269/297 (90%)	1.20	56 (20%) 2 4	116, 163, 212, 229	0
1	D	268/297 (90%)	1.82	78 (29%) 1 2	110, 168, 216, 236	0
1	E	269/297 (90%)	0.88	45 (16%) 4 7	104, 145, 213, 231	0
All	All	1348/1485 (90%)	1.10	238 (17%) 4 6	104, 158, 213, 261	0

All (238) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	49	GLY	16.5
1	D	170	GLN	16.1
1	D	265	THR	14.8
1	D	145	GLU	14.5
1	B	46	GLU	14.5
1	C	106	ALA	14.2
1	B	47	GLN	13.7
1	B	48	LEU	13.6
1	D	146	ARG	12.2
1	D	167	GLU	12.2
1	C	105	PRO	11.8
1	A	47	GLN	11.4
1	D	166	ASN	10.9
1	C	109	ASP	10.6
1	D	236	TYR	10.6
1	D	112	ARG	10.5
1	D	47	GLN	10.2
1	E	170	GLN	9.6
1	D	173	GLU	9.6
1	D	174	ALA	9.4

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Mol	Chain	Res	Type	RSRZ
1	C	101	ARG	9.2
1	D	149	GLU	9.1
1	A	48	LEU	9.0
1	D	113	ARG	8.6
1	C	204	THR	8.3
1	C	25	ILE	8.0
1	E	174	ALA	7.7
1	D	110	ALA	7.4
1	E	173	GLU	7.2
1	D	233	ASP	7.1
1	D	232	GLY	7.1
1	D	46	GLU	6.8
1	B	50	ILE	6.8
1	A	174	ALA	6.6
1	D	117	TYR	6.5
1	C	211	TYR	6.4
1	C	209	PHE	6.4
1	D	235	HIS	6.2
1	A	175	GLY	6.2
1	B	44	TRP	6.1
1	B	216	GLN	6.1
1	D	176	LYS	6.0
1	D	42	TYR	6.0
1	E	47	GLN	5.9
1	A	145	GLU	5.9
1	D	163	LEU	5.9
1	D	154	SER	5.8
1	B	51	HIS	5.7
1	C	285	ASP	5.7
1	B	211	TYR	5.6
1	D	266	ALA	5.6
1	E	175	GLY	5.5
1	A	281	ARG	5.4
1	A	144	GLU	5.4
1	C	24	ILE	5.4
1	B	260	GLU	5.2
1	C	257	GLU	5.2
1	E	42	TYR	5.2
1	C	29	LEU	5.1
1	D	107	GLU	5.1
1	D	116	SER	5.0
1	C	107	GLU	5.0

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Mol	Chain	Res	Type	RSRZ
1	C	184	LEU	4.9
1	D	32	VAL	4.7
1	C	27	ARG	4.7
1	E	172	ARG	4.6
1	B	70	PHE	4.6
1	D	277	ASN	4.5
1	B	45	TYR	4.5
1	B	257	GLU	4.4
1	E	26	PHE	4.4
1	D	106	ALA	4.4
1	A	176	LYS	4.4
1	E	154	SER	4.3
1	A	70	PHE	4.3
1	D	132	THR	4.3
1	B	288	GLY	4.2
1	B	287	THR	4.1
1	B	73	SER	4.1
1	E	266	ALA	4.1
1	D	49	GLY	4.1
1	E	262	ALA	4.0
1	D	187	ASN	4.0
1	D	109	ASP	4.0
1	D	264	GLY	3.9
1	C	253	ASP	3.9
1	E	152	ALA	3.8
1	E	169	GLY	3.8
1	A	102	ASN	3.8
1	D	143	PRO	3.8
1	D	159	ARG	3.8
1	E	53	THR	3.7
1	A	211	TYR	3.7
1	B	77	SER	3.7
1	A	173	GLU	3.7
1	E	290	HIS	3.6
1	D	224	THR	3.6
1	D	171	LEU	3.6
1	D	184	LEU	3.5
1	C	67	PHE	3.5
1	A	103	ILE	3.5
1	D	34	MET	3.5
1	D	130	ARG	3.5
1	C	232	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	239	PRO	3.5
1	D	142	LEU	3.4
1	E	253	ASP	3.4
1	C	180	ILE	3.3
1	D	175	GLY	3.3
1	C	205	THR	3.3
1	E	45	TYR	3.3
1	D	153	SER	3.3
1	C	137	ASP	3.2
1	A	46	GLU	3.2
1	E	52	LEU	3.2
1	D	268	ASN	3.2
1	B	42	TYR	3.2
1	E	252	TRP	3.2
1	D	211	TYR	3.2
1	D	131	LYS	3.2
1	C	210	ALA	3.2
1	D	43	GLN	3.1
1	B	285	ASP	3.1
1	E	249	PHE	3.1
1	A	204	THR	3.1
1	B	52	LEU	3.1
1	E	48	LEU	3.1
1	B	209	PHE	3.1
1	A	148	THR	3.1
1	C	108	HIS	3.1
1	B	215	LEU	3.0
1	B	212	THR	3.0
1	C	187	ASN	3.0
1	E	176	LYS	3.0
1	C	288	GLY	3.0
1	D	193	ALA	3.0
1	D	48	LEU	3.0
1	A	233	ASP	3.0
1	C	256	ALA	2.9
1	E	259	LEU	2.9
1	C	212	THR	2.9
1	A	106	ALA	2.9
1	C	286	MET	2.9
1	C	260	GLU	2.9
1	A	44	TRP	2.9
1	B	193	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	183	GLY	2.8
1	B	219	VAL	2.8
1	D	183	GLY	2.8
1	E	235	HIS	2.8
1	B	197	GLY	2.8
1	D	280	GLU	2.8
1	B	93	GLU	2.7
1	E	232	GLY	2.7
1	C	185	MET	2.7
1	A	257	GLU	2.7
1	D	179	ASP	2.7
1	D	186	ASP	2.6
1	B	32	VAL	2.6
1	E	207	VAL	2.6
1	C	206	PRO	2.6
1	C	181	THR	2.6
1	C	72	ASN	2.6
1	D	168	ILE	2.6
1	A	278	THR	2.6
1	C	30	LEU	2.6
1	E	219	VAL	2.5
1	D	28	LEU	2.5
1	D	243	VAL	2.5
1	C	26	PHE	2.5
1	E	220	TYR	2.5
1	C	46	GLU	2.5
1	D	223	CYS	2.5
1	C	208	PRO	2.5
1	D	180	ILE	2.5
1	D	37	ILE	2.5
1	A	205	THR	2.5
1	C	215	LEU	2.5
1	E	171	LEU	2.5
1	D	25	ILE	2.5
1	E	83	ARG	2.5
1	C	216	GLN	2.4
1	D	31	ASN	2.4
1	E	263	PHE	2.4
1	D	141	LEU	2.4
1	D	70	PHE	2.4
1	C	276	CYS	2.4
1	D	82	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	144	GLU	2.4
1	D	267	ALA	2.4
1	E	275	MET	2.4
1	E	70	PHE	2.4
1	C	230	LEU	2.4
1	D	50	ILE	2.4
1	E	166	ASN	2.3
1	C	96	LEU	2.3
1	E	39	ILE	2.3
1	C	94	ARG	2.3
1	E	64	ILE	2.3
1	C	188	LYS	2.3
1	D	263	PHE	2.3
1	C	51	HIS	2.3
1	D	29	LEU	2.3
1	E	177	LEU	2.3
1	B	66	ILE	2.3
1	D	206	PRO	2.3
1	D	104	LEU	2.3
1	D	210	ALA	2.3
1	C	289	GLN	2.2
1	B	101	ARG	2.2
1	E	109	ASP	2.2
1	C	254	SER	2.2
1	E	153	SER	2.2
1	C	233	ASP	2.2
1	B	119	VAL	2.1
1	A	209	PHE	2.1
1	A	235	HIS	2.1
1	D	164	ALA	2.1
1	C	214	ILE	2.1
1	D	150	ILE	2.1
1	C	134	PRO	2.1
1	D	30	LEU	2.1
1	A	92	ALA	2.1
1	E	29	LEU	2.1
1	E	66	ILE	2.1
1	D	225	LEU	2.1
1	E	65	ALA	2.1
1	C	48	LEU	2.1
1	E	67	PHE	2.0
1	B	225	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	99	GLN	2.0
1	B	69	GLY	2.0
1	C	261	ASP	2.0
1	C	73	SER	2.0
1	E	254	SER	2.0
1	C	213	LEU	2.0
1	A	262	ALA	2.0
1	E	233	ASP	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 ⓘ

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(Å ²)	Q<0.9
2	ZN	C	305	1/1	0.29	0.17	260,260,260,260	0
2	ZN	D	303	1/1	0.88	0.17	230,230,230,230	0
3	CL	B	303	1/1	0.88	0.09	182,182,182,182	0
3	CL	C	304	1/1	0.89	0.19	162,162,162,162	0
2	ZN	D	302	1/1	0.91	0.11	169,169,169,169	0
2	ZN	C	302	1/1	0.94	0.15	148,148,148,148	0
3	CL	B	304	1/1	0.95	0.09	151,151,151,151	0
2	ZN	A	301	1/1	0.95	0.14	117,117,117,117	0
3	CL	A	305	1/1	0.96	0.15	103,103,103,103	0
2	ZN	B	301	1/1	0.96	0.06	198,198,198,198	0
3	CL	D	305	1/1	0.96	0.04	150,150,150,150	0
3	CL	D	306	1/1	0.96	0.15	191,191,191,191	0
3	CL	A	304	1/1	0.97	0.11	141,141,141,141	0
3	CL	C	303	1/1	0.98	0.06	115,115,115,115	0
2	ZN	A	303	1/1	0.98	0.07	132,132,132,132	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	B	302	1/1	0.98	0.04	146,146,146,146	0
2	ZN	C	306	1/1	0.98	0.13	112,112,112,112	0
3	CL	E	302	1/1	0.98	0.05	116,116,116,116	0
2	ZN	A	302	1/1	0.99	0.03	120,120,120,120	0
3	CL	D	304	1/1	0.99	0.05	108,108,108,108	0
2	ZN	B	305	1/1	0.99	0.04	132,132,132,132	0
2	ZN	E	301	1/1	0.99	0.02	120,120,120,120	0
2	ZN	C	301	1/1	0.99	0.07	124,124,124,124	0
2	ZN	D	301	1/1	1.00	0.03	167,167,167,167	0

6.5 Other polymers [i](#)

There are no such residues in this entry.