



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 11:10 AM UTC

PDB ID : 6IV3 / pdb_00006iv3
Title : Crystal structure of a bacterial Bestrophin homolog from *Klebsiella pneumoniae* with a mutation W252A
Authors : Kittredge, A.; Chen, S.; Yang, T.
Deposited on : 2018-12-02
Resolution : 2.52 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

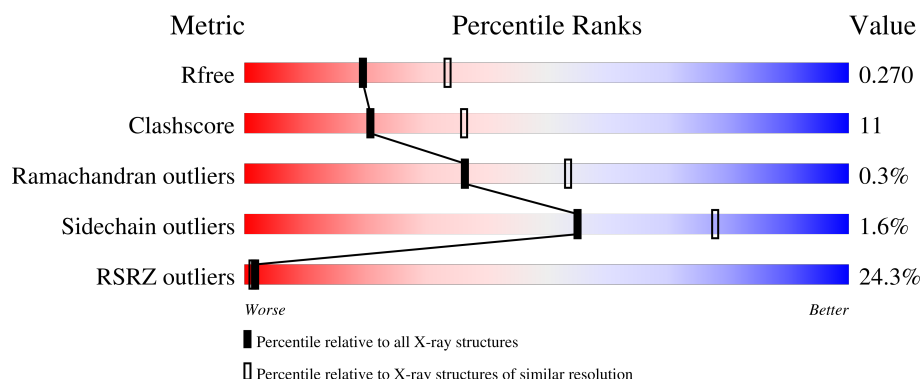
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.52 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	7383 (2.54-2.50)
Clashscore	190562	8079 (2.54-2.50)
Ramachandran outliers	187476	7944 (2.54-2.50)
Sidechain outliers	187428	7946 (2.54-2.50)
RSRZ outliers	180081	7387 (2.54-2.50)

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	296	17% 71% 22% 7%
1	B	296	18% 77% 15% 8%
1	C	296	23% 73% 18% 8%
1	D	296	34% 69% 21% 9%
1	E	296	19% 68% 20% 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZN	D	304	-	-	-	X
2	ZN	E	303	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10752 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 homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2160	1398	367	386	9			
1	B	272	Total	C	N	O	S	0	0	0
			2159	1398	366	386	9			
1	C	271	Total	C	N	O	S	0	0	0
			2112	1362	361	380	9			
1	D	268	Total	C	N	O	S	0	0	0
			2109	1361	358	381	9			
1	E	270	Total	C	N	O	S	0	0	0
			2117	1369	358	381	9			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ASN	-	expression tag	UNP W1ELP7
A	0	ALA	-	expression tag	UNP W1ELP7
A	252	ALA	TRP	engineered mutation	UNP W1ELP7
B	-1	ASN	-	expression tag	UNP W1ELP7
B	0	ALA	-	expression tag	UNP W1ELP7
B	252	ALA	TRP	engineered mutation	UNP W1ELP7
C	-1	ASN	-	expression tag	UNP W1ELP7
C	0	ALA	-	expression tag	UNP W1ELP7
C	252	ALA	TRP	engineered mutation	UNP W1ELP7
D	-1	ASN	-	expression tag	UNP W1ELP7
D	0	ALA	-	expression tag	UNP W1ELP7
D	252	ALA	TRP	engineered mutation	UNP W1ELP7
E	-1	ASN	-	expression tag	UNP W1ELP7
E	0	ALA	-	expression tag	UNP W1ELP7
E	252	ALA	TRP	engineered mutation	UNP W1ELP7

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

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

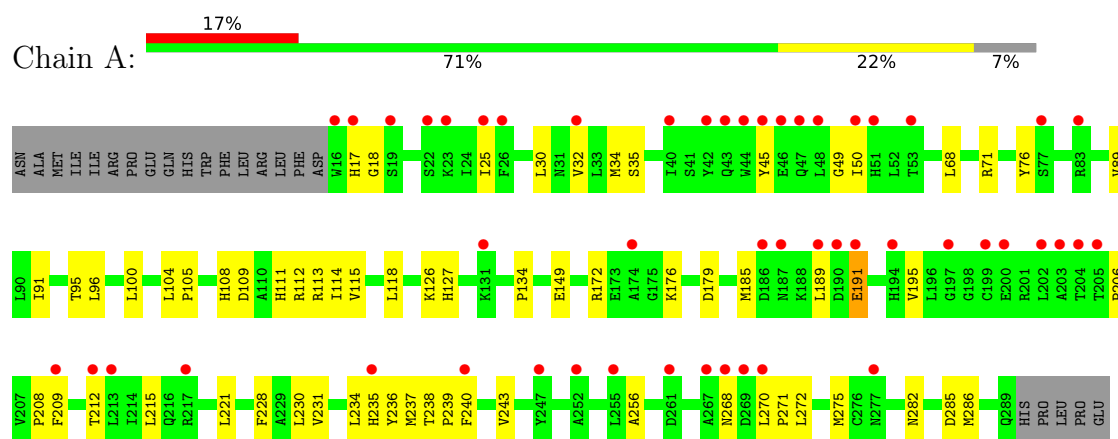
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	26	Total 26	O 26	0	0
3	B	13	Total 13	O 13	0	0
3	C	18	Total 18	O 18	0	0
3	D	8	Total 8	O 8	0	0
3	E	15	Total 15	O 15	0	0

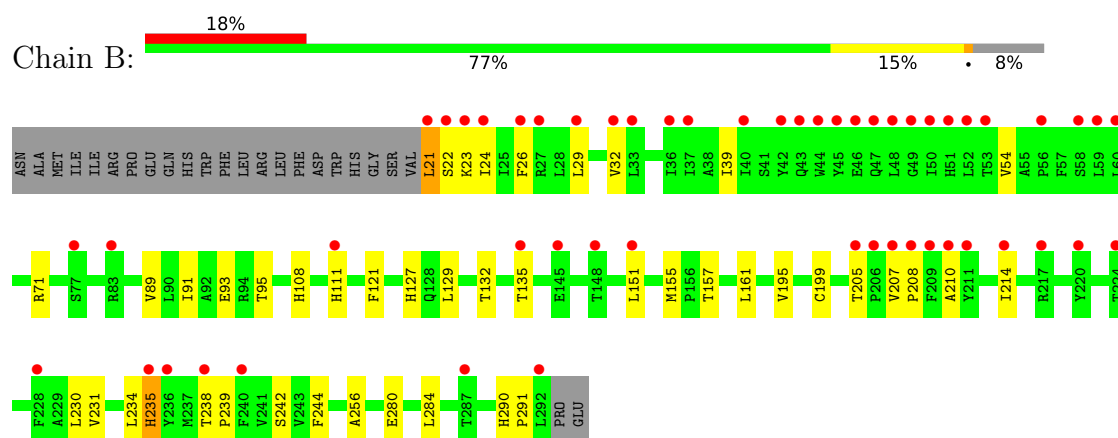
3 Residue-property plots [i](#)

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.

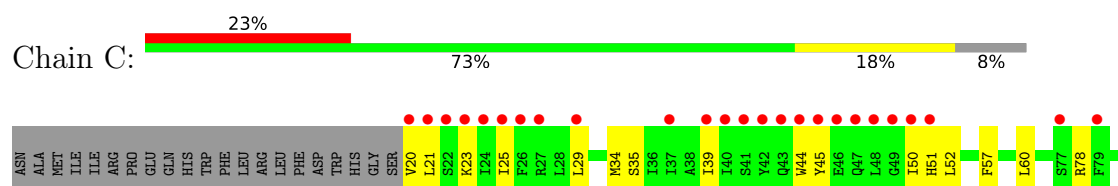
• Molecule 1: Bestrophin homolog

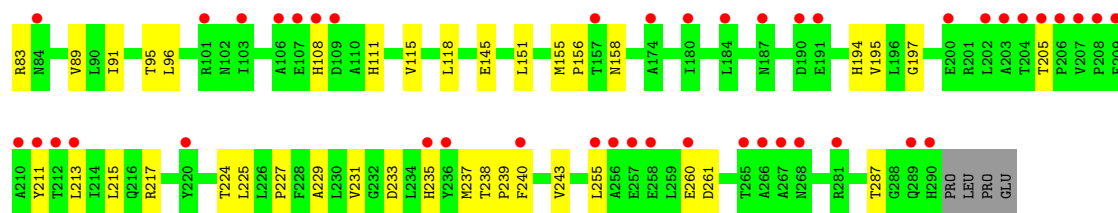


• Molecule 1: Bestrophin homolog

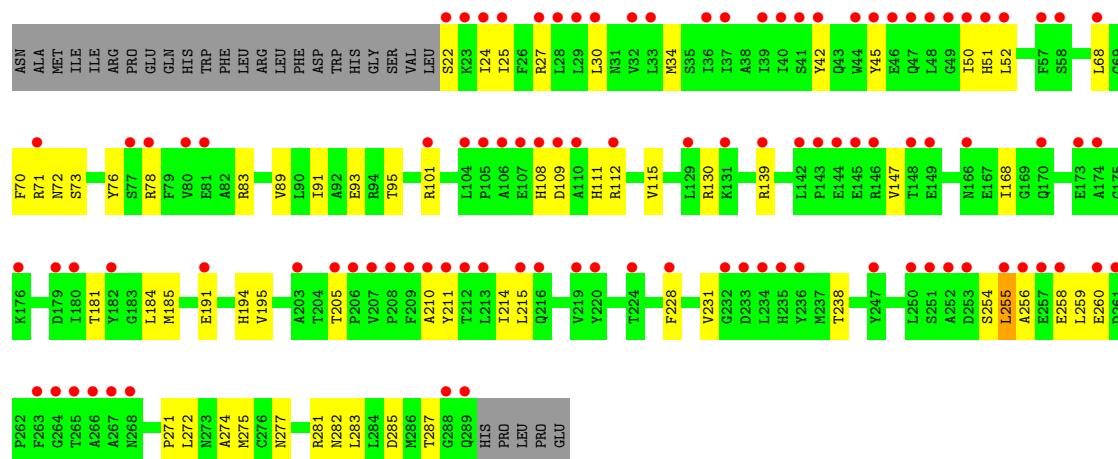


• Molecule 1: Bestrophin homolog

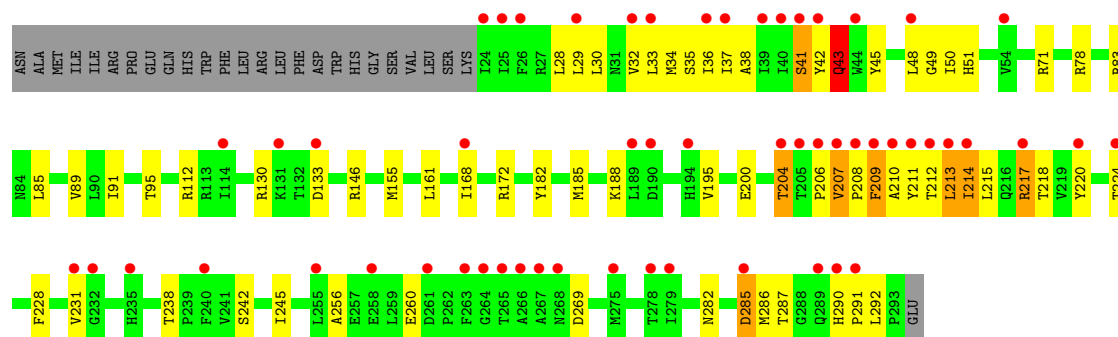




• Molecule 1: Bestrophin homolog



• Molecule 1: Bestrophin homolog



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	113.46Å 159.61Å 162.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.11 – 2.52 66.11 – 2.52	Depositor EDS
% Data completeness (in resolution range)	98.7 (66.11-2.52) 98.6 (66.11-2.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.234 , 0.265 0.242 , 0.270	Depositor DCC
R_{free} test set	2007 reflections (1.99%)	wwPDB-VP
Wilson B-factor (Å ²)	62.4	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 36.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10752	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.50	0/2207	0.71	5/3006 (0.2%)
1	B	0.38	0/2205	0.59	0/3001
1	C	0.43	0/2153	0.61	0/2932
1	D	0.46	0/2152	0.62	0/2930
1	E	0.56	2/2161 (0.1%)	0.74	1/2945 (0.0%)
All	All	0.47	2/10878 (0.0%)	0.66	6/14814 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	218	THR	C-O	-6.06	1.17	1.24
1	E	215	LEU	C-O	-5.35	1.17	1.24

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	209	PHE	N-CA-C	-6.90	104.86	113.20
1	A	189	LEU	CA-C-N	-6.79	110.08	122.38
1	A	189	LEU	C-N-CA	-6.79	110.08	122.38
1	A	191	GLU	N-CA-C	-6.17	104.63	111.36
1	A	17	HIS	CA-C-N	5.16	131.52	121.41
1	A	17	HIS	C-N-CA	5.16	131.52	121.41

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	2160	0	2191	62	0
1	B	2159	0	2212	41	0
1	C	2112	0	2152	44	0
1	D	2109	0	2143	57	0
1	E	2117	0	2155	70	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
2	C	3	0	0	0	0
2	D	4	0	0	0	0
2	E	3	0	0	0	0
3	A	26	0	0	0	0
3	B	13	0	0	0	0
3	C	18	0	0	2	0
3	D	8	0	0	1	0
3	E	15	0	0	2	0
All	All	10752	0	10853	239	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 11.

All (239) 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:72:ASN:OD1	1:D:255:LEU:HD23	1.55	1.05
1:E:211:TYR:O	1:E:214:ILE:HG23	1.62	1.00
1:E:285:ASP:HA	1:E:290:HIS:CB	2.05	0.87
1:B:231:VAL:HA	1:B:238:THR:HG21	1.59	0.84
1:B:234:LEU:HB2	1:B:238:THR:HG22	1.62	0.82
1:C:45:TYR:O	1:C:50:ILE:HG12	1.80	0.81
1:D:71:ARG:CZ	1:D:256:ALA:HB3	2.10	0.80
1:E:211:TYR:O	1:E:214:ILE:CG2	2.30	0.79
1:A:268:ASN:ND2	1:E:210:ALA:HB2	1.97	0.78
1:C:155:MET:HG2	1:C:158:ASN:HB2	1.66	0.76
1:D:71:ARG:CZ	1:D:256:ALA:CB	2.64	0.75
1:D:101:ARG:HH21	1:D:287:THR:HA	1.51	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:ARG:HG2	1:D:205:THR:HG21	1.69	0.74
1:C:91:ILE:HD13	1:D:194:HIS:HA	1.69	0.74
1:C:21:LEU:C	1:C:21:LEU:HD13	2.13	0.74
1:D:71:ARG:NH2	1:D:256:ALA:HB3	2.03	0.74
1:D:83:ARG:NH2	1:D:83:ARG:O	2.20	0.73
1:D:72:ASN:OD1	1:D:255:LEU:CD2	2.34	0.73
1:E:168:ILE:HD12	1:E:185:MET:HE3	1.71	0.72
1:B:205:THR:HG21	1:D:83:ARG:HE	1.56	0.70
1:E:34:MET:HE3	1:E:228:PHE:HE1	1.54	0.70
1:A:268:ASN:ND2	1:E:210:ALA:CB	2.54	0.69
1:A:230:LEU:HB3	1:A:234:LEU:HD12	1.74	0.68
1:B:135:THR:HG22	1:B:151:LEU:HD11	1.76	0.68
1:E:188:LYS:NZ	3:E:401:HOH:O	2.27	0.68
1:D:71:ARG:NE	1:D:256:ALA:CB	2.56	0.67
1:E:95:THR:HG21	1:E:188:LYS:HE2	1.76	0.67
1:D:22:SER:O	1:D:25:ILE:HG22	1.96	0.66
1:C:21:LEU:HD13	1:C:21:LEU:O	1.96	0.65
1:D:115:VAL:HG21	1:D:287:THR:HG21	1.78	0.65
1:B:108:HIS:HA	1:B:111:HIS:CG	2.31	0.65
1:D:211:TYR:O	1:D:215:LEU:HB2	1.97	0.64
1:A:272:LEU:HD23	1:A:275:MET:HE3	1.80	0.64
1:A:91:ILE:O	1:A:95:THR:HG23	1.97	0.63
1:A:270:LEU:HB3	1:A:275:MET:HE1	1.79	0.63
1:A:285:ASP:CB	1:E:155:MET:HE1	2.28	0.63
1:B:205:THR:HG21	1:D:83:ARG:NE	2.13	0.63
1:C:45:TYR:CB	1:C:50:ILE:HD11	2.28	0.63
1:B:91:ILE:O	1:B:95:THR:HG23	1.99	0.62
1:A:285:ASP:HB3	1:E:155:MET:HE1	1.80	0.62
1:B:111:HIS:CE1	1:B:290:HIS:HE2	2.17	0.62
1:B:111:HIS:CE1	1:B:290:HIS:NE2	2.68	0.61
1:E:211:TYR:HA	1:E:214:ILE:CG2	2.30	0.61
1:E:285:ASP:CA	1:E:290:HIS:CB	2.78	0.61
1:C:194:HIS:HA	1:E:91:ILE:HD13	1.81	0.61
1:E:282:ASN:O	1:E:286:MET:HB2	2.01	0.61
1:B:155:MET:HE1	1:D:285:ASP:HB3	1.82	0.60
1:E:207:VAL:O	1:E:208:PRO:C	2.43	0.60
1:C:205:THR:OG1	1:E:83:ARG:NH1	2.35	0.59
1:B:155:MET:HE3	1:D:282:ASN:HA	1.85	0.58
1:E:214:ILE:C	1:E:214:ILE:HD13	2.28	0.58
1:C:20:VAL:HG13	1:C:23:LYS:H	1.69	0.58
1:A:109:ASP:OD1	1:A:112:ARG:NH2	2.37	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:45:TYR:HA	1:E:48:LEU:HD22	1.86	0.57
1:A:206:PRO:HG3	1:A:209:PHE:CZ	2.39	0.57
1:C:213:LEU:HD11	1:C:217:ARG:NH1	2.19	0.57
1:D:24:ILE:HD11	1:D:254:SER:HB3	1.85	0.57
1:B:111:HIS:HE1	1:B:290:HIS:CE1	2.23	0.57
1:E:34:MET:HE3	1:E:228:PHE:CE1	2.38	0.57
1:D:109:ASP:HA	1:D:112:ARG:HE	1.69	0.56
1:A:68:LEU:HD23	1:A:215:LEU:HD13	1.88	0.56
1:C:237:MET:HE2	1:D:52:LEU:HD21	1.86	0.56
1:D:91:ILE:O	1:D:95:THR:HG23	2.06	0.56
1:B:111:HIS:CE1	1:B:290:HIS:CE1	2.94	0.56
1:E:37:ILE:O	1:E:41:SER:HB2	2.06	0.55
1:C:78:ARG:NH1	1:C:260:GLU:O	2.37	0.55
1:D:27:ARG:HD3	1:D:30:LEU:HD23	1.89	0.55
1:B:280:GLU:HG3	1:B:284:LEU:HD12	1.89	0.55
1:A:235:HIS:HB2	1:E:51:HIS:CD2	2.42	0.55
1:D:256:ALA:O	1:D:260:GLU:HG2	2.06	0.55
1:C:145:GLU:CD	1:C:145:GLU:H	2.15	0.54
1:E:71:ARG:HD3	1:E:256:ALA:HB1	1.89	0.54
1:D:78:ARG:NH1	1:D:205:THR:O	2.41	0.54
1:A:235:HIS:HB3	1:E:49:GLY:O	2.08	0.54
1:E:89:VAL:CG2	1:E:195:VAL:HG11	2.37	0.54
1:C:57:PHE:HE2	1:C:225:LEU:HB3	1.73	0.54
1:A:282:ASN:HA	1:E:155:MET:HE3	1.89	0.53
1:D:93:GLU:HB3	1:D:283:LEU:HD11	1.89	0.53
1:A:96:LEU:CD1	1:A:185:MET:HB3	2.39	0.53
1:E:45:TYR:O	1:E:48:LEU:HB2	2.09	0.53
1:E:71:ARG:NH2	1:E:260:GLU:OE1	2.35	0.53
1:E:38:ALA:O	1:E:42:TYR:HB2	2.09	0.53
1:B:21:LEU:HD13	1:B:24:ILE:HG23	1.90	0.52
1:A:237:MET:HG2	1:E:50:ILE:HG23	1.91	0.52
1:D:89:VAL:HG23	1:D:195:VAL:HG11	1.92	0.52
1:A:100:LEU:HG	1:A:185:MET:HE1	1.90	0.52
1:B:108:HIS:O	1:B:111:HIS:HB2	2.10	0.52
1:B:210:ALA:HA	1:B:214:ILE:HD13	1.90	0.52
1:D:89:VAL:CG2	1:D:195:VAL:HG11	2.40	0.52
1:E:78:ARG:CD	1:E:207:VAL:HG12	2.40	0.52
1:A:76:TYR:HE1	1:E:207:VAL:CG2	2.22	0.51
1:A:270:LEU:HB3	1:A:275:MET:CE	2.40	0.51
1:B:127:HIS:ND1	1:B:132:THR:OG1	2.38	0.51
1:E:290:HIS:N	1:E:291:PRO:CD	2.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:235:HIS:CD2	1:D:51:HIS:CE1	2.98	0.51
1:E:78:ARG:NH2	1:E:206:PRO:HA	2.26	0.51
1:D:271:PRO:HB2	1:D:274:ALA:HB3	1.92	0.50
1:D:108:HIS:HA	1:D:111:HIS:CG	2.46	0.50
1:C:235:HIS:CD2	1:D:51:HIS:HE1	2.29	0.50
1:E:42:TYR:O	1:E:43:GLN:HB2	2.11	0.50
1:E:78:ARG:HD2	1:E:207:VAL:HG12	1.94	0.50
1:E:209:PHE:HA	1:E:212:THR:OG1	2.12	0.50
1:A:76:TYR:HE1	1:E:207:VAL:HG23	1.76	0.49
1:C:21:LEU:C	1:C:21:LEU:CD1	2.83	0.49
1:A:32:VAL:HG23	1:A:243:VAL:HG21	1.94	0.49
1:A:34:MET:HE3	1:A:228:PHE:CE1	2.47	0.49
1:D:271:PRO:HD2	1:D:275:MET:HE3	1.92	0.49
1:B:238:THR:O	1:B:242:SER:CB	2.61	0.49
1:A:104:LEU:O	1:A:111:HIS:NE2	2.44	0.49
1:A:71:ARG:HH11	1:A:212:THR:HG22	1.78	0.49
1:A:76:TYR:CE1	1:E:207:VAL:CG2	2.96	0.49
1:E:200:GLU:O	1:E:204:THR:OG1	2.20	0.49
1:C:197:GLY:HA3	1:E:91:ILE:HD11	1.95	0.49
1:A:109:ASP:HB3	1:A:113:ARG:HH12	1.77	0.48
1:A:271:PRO:HD2	1:A:275:MET:HE2	1.95	0.48
1:D:30:LEU:O	1:D:34:MET:HG2	2.12	0.48
1:D:108:HIS:HA	1:D:111:HIS:ND1	2.28	0.48
1:D:181:THR:HA	1:D:184:LEU:HD12	1.95	0.48
1:C:213:LEU:HD11	1:C:217:ARG:CZ	2.43	0.48
1:E:155:MET:HB2	1:E:155:MET:HE2	1.75	0.48
1:D:101:ARG:NH1	1:D:108:HIS:CE1	2.81	0.48
1:E:211:TYR:C	1:E:214:ILE:CG2	2.86	0.48
1:A:76:TYR:OH	1:E:207:VAL:HB	2.13	0.48
1:B:29:LEU:HA	1:B:32:VAL:HG23	1.94	0.48
1:A:45:TYR:CD1	1:A:50:ILE:HD11	2.49	0.48
1:C:89:VAL:CG2	1:C:195:VAL:HG11	2.44	0.48
1:E:89:VAL:HG23	1:E:195:VAL:HG11	1.95	0.48
1:A:271:PRO:O	1:A:275:MET:HG3	2.14	0.47
1:C:231:VAL:HA	1:C:238:THR:OG1	2.14	0.47
1:D:231:VAL:HA	1:D:238:THR:OG1	2.15	0.47
1:C:240:PHE:O	1:C:243:VAL:HG12	2.15	0.47
1:A:231:VAL:HA	1:A:238:THR:OG1	2.15	0.47
1:A:268:ASN:CG	1:E:210:ALA:HB2	2.40	0.47
1:A:271:PRO:CD	1:A:275:MET:HE2	2.44	0.47
1:B:135:THR:HG22	1:B:151:LEU:CD1	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:255:LEU:O	1:D:255:LEU:HG	2.11	0.46
1:A:282:ASN:O	1:A:286:MET:HG3	2.14	0.46
1:D:70:PHE:O	1:D:73:SER:HB3	2.14	0.46
1:B:155:MET:HE1	1:D:285:ASP:CB	2.46	0.46
1:D:277:ASN:O	1:D:281:ARG:HG3	2.15	0.46
1:B:238:THR:O	1:B:242:SER:HB2	2.15	0.46
1:C:211:TYR:O	1:C:215:LEU:HB2	2.16	0.46
1:D:68:LEU:HD23	1:D:215:LEU:HD13	1.97	0.46
1:A:32:VAL:CG2	1:A:243:VAL:HG21	2.46	0.46
1:C:260:GLU:HG3	1:C:261:ASP:N	2.31	0.46
1:A:49:GLY:O	1:B:235:HIS:HB2	2.15	0.46
1:A:237:MET:HE1	1:A:240:PHE:HB2	1.98	0.46
1:B:129:LEU:HD21	1:B:199:CYS:HB3	1.98	0.46
1:B:161:LEU:HD23	1:B:161:LEU:HA	1.80	0.45
1:D:71:ARG:NH2	1:D:256:ALA:CB	2.76	0.45
1:C:44:TRP:CD1	1:C:44:TRP:N	2.84	0.45
1:C:35:SER:HA	1:C:227:PRO:HB2	1.99	0.45
1:D:101:ARG:HD3	3:D:406:HOH:O	2.17	0.45
1:C:57:PHE:CE2	1:C:225:LEU:HB3	2.49	0.45
1:E:211:TYR:HA	1:E:214:ILE:HG21	1.99	0.45
1:C:89:VAL:HG23	1:C:195:VAL:HG11	1.99	0.45
1:E:32:VAL:O	1:E:35:SER:HB2	2.16	0.45
1:E:146:ARG:HA	1:E:146:ARG:NE	2.33	0.44
1:B:121:PHE:CE1	1:B:157:THR:HG22	2.52	0.44
1:B:238:THR:N	1:B:239:PRO:HD2	2.32	0.44
1:C:34:MET:HG3	1:C:224:THR:HG23	2.00	0.44
1:A:89:VAL:CG2	1:A:195:VAL:HG11	2.48	0.44
1:A:105:PRO:O	1:A:108:HIS:CE1	2.70	0.44
1:C:91:ILE:O	1:C:95:THR:HG23	2.18	0.44
1:E:231:VAL:HA	1:E:238:THR:OG1	2.18	0.44
1:A:96:LEU:HD11	1:A:185:MET:HB3	1.99	0.44
1:C:156:PRO:HD2	3:C:405:HOH:O	2.18	0.44
1:B:39:ILE:HD11	1:B:239:PRO:HD3	2.00	0.43
1:C:115:VAL:HG21	1:C:287:THR:HG21	1.98	0.43
1:C:233:ASP:CG	3:C:401:HOH:O	2.60	0.43
1:A:104:LEU:HD21	1:A:176:LYS:O	2.18	0.43
1:D:215:LEU:HD23	1:D:215:LEU:HA	1.92	0.43
1:E:130:ARG:HH21	1:E:269:ASP:HB2	1.83	0.43
1:E:172:ARG:NH2	3:E:402:HOH:O	2.29	0.43
1:D:34:MET:HE3	1:D:228:PHE:HE2	1.83	0.43
1:A:114:ILE:O	1:A:118:LEU:HG	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:VAL:HG23	1:D:76:TYR:CE1	2.53	0.43
1:D:210:ALA:O	1:D:214:ILE:HD12	2.18	0.43
1:E:210:ALA:O	1:E:214:ILE:HG22	2.19	0.43
1:E:85:LEU:HD23	1:E:85:LEU:HA	1.90	0.43
1:C:25:ILE:O	1:C:29:LEU:HD23	2.19	0.43
1:A:96:LEU:HD23	1:A:118:LEU:HD11	2.00	0.43
1:C:45:TYR:C	1:C:50:ILE:HG12	2.44	0.43
1:E:238:THR:O	1:E:242:SER:HB2	2.19	0.43
1:B:89:VAL:CG2	1:B:195:VAL:HG11	2.49	0.42
1:A:71:ARG:CZ	1:A:256:ALA:HB2	2.49	0.42
1:B:108:HIS:HA	1:B:111:HIS:ND1	2.34	0.42
1:D:168:ILE:HD12	1:D:185:MET:HE2	1.99	0.42
1:E:220:TYR:O	1:E:224:THR:HG23	2.18	0.42
1:A:237:MET:HE3	1:A:237:MET:O	2.19	0.42
1:E:95:THR:CG2	1:E:188:LYS:HE2	2.47	0.42
1:A:172:ARG:NH2	1:A:179:ASP:OD1	2.46	0.42
1:D:42:TYR:O	1:D:45:TYR:HB2	2.20	0.42
1:A:127:HIS:CD2	1:A:134:PRO:HA	2.55	0.42
1:B:71:ARG:HD2	1:B:256:ALA:HB1	2.01	0.42
1:E:211:TYR:C	1:E:214:ILE:HG22	2.45	0.42
1:D:83:ARG:HA	1:D:83:ARG:HD2	1.82	0.42
1:A:268:ASN:HD21	1:E:210:ALA:CB	2.31	0.42
1:B:208:PRO:HG2	1:D:259:LEU:HD21	2.01	0.42
1:E:161:LEU:HA	1:E:161:LEU:HD23	1.75	0.42
1:E:211:TYR:CA	1:E:214:ILE:CG2	2.98	0.42
1:C:151:LEU:HD23	1:C:151:LEU:HA	1.84	0.42
1:A:45:TYR:CE1	1:A:50:ILE:HD11	2.54	0.41
1:A:221:LEU:HB3	1:B:244:PHE:CZ	2.55	0.41
1:A:268:ASN:ND2	1:E:210:ALA:HB1	2.32	0.41
1:B:23:LYS:O	1:B:26:PHE:HB3	2.21	0.41
1:A:230:LEU:HB3	1:A:234:LEU:CD1	2.48	0.41
1:C:39:ILE:HD11	1:C:239:PRO:HD3	2.01	0.41
1:C:51:HIS:C	1:C:52:LEU:HD12	2.45	0.41
1:C:96:LEU:HD23	1:C:118:LEU:HD21	2.02	0.41
1:D:34:MET:HE3	1:D:228:PHE:CE2	2.55	0.41
1:D:108:HIS:O	1:D:111:HIS:HB2	2.20	0.41
1:E:29:LEU:O	1:E:32:VAL:HG22	2.20	0.41
1:E:33:LEU:O	1:E:36:ILE:HG12	2.20	0.41
1:A:71:ARG:NH1	1:A:212:THR:HG22	2.35	0.41
1:A:100:LEU:HG	1:A:185:MET:CE	2.50	0.41
1:B:54:VAL:HG22	1:B:230:LEU:HD23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:255:LEU:HD12	1:C:255:LEU:HA	1.85	0.41
1:D:139:ARG:HG2	1:D:147:VAL:HG11	2.02	0.41
1:E:213:LEU:HD22	1:E:213:LEU:HA	1.63	0.41
1:B:22:SER:OG	1:B:23:LYS:HD2	2.21	0.41
1:C:108:HIS:HA	1:C:111:HIS:ND1	2.36	0.41
1:A:111:HIS:O	1:A:115:VAL:HG23	2.21	0.41
1:A:149:GLU:HG2	1:B:291:PRO:HG3	2.01	0.41
1:E:168:ILE:HG22	1:E:182:TYR:CD1	2.55	0.41
1:C:60:LEU:HD13	1:E:245:ILE:HA	2.02	0.41
1:D:50:ILE:HG13	1:D:50:ILE:O	2.21	0.41
1:E:217:ARG:HA	1:E:217:ARG:HD2	1.78	0.41
1:A:126:LYS:HD2	1:A:126:LYS:C	2.45	0.41
1:A:236:TYR:HB2	1:E:50:ILE:HD13	2.03	0.41
1:B:89:VAL:O	1:B:93:GLU:HG3	2.20	0.41
1:A:30:LEU:O	1:A:34:MET:HG2	2.21	0.41
1:A:215:LEU:HD23	1:A:215:LEU:HA	1.96	0.41
1:C:52:LEU:HB3	1:C:229:ALA:HA	2.01	0.41
1:B:210:ALA:O	1:B:214:ILE:HB	2.22	0.40
1:E:30:LEU:O	1:E:34:MET:HG2	2.21	0.40
1:D:130:ARG:NH2	1:D:272:LEU:HB2	2.37	0.40
1:A:35:SER:OG	1:A:239:PRO:HB3	2.21	0.40
1:A:71:ARG:NE	1:A:256:ALA:HB2	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	272/296 (92%)	264 (97%)	6 (2%)	2 (1%)	18	32
1	B	270/296 (91%)	262 (97%)	7 (3%)	1 (0%)	30	47
1	C	269/296 (91%)	265 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	266/296 (90%)	257 (97%)	9 (3%)	0	100	100
1	E	268/296 (90%)	264 (98%)	3 (1%)	1 (0%)	30	47
All	All	1345/1480 (91%)	1312 (98%)	29 (2%)	4 (0%)	36	53

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	GLY
1	E	43	GLN
1	B	235	HIS
1	A	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	233/258 (90%)	231 (99%)	2 (1%)	70	86
1	B	236/258 (92%)	235 (100%)	1 (0%)	84	92
1	C	227/258 (88%)	227 (100%)	0	100	100
1	D	228/258 (88%)	225 (99%)	3 (1%)	61	81
1	E	229/258 (89%)	216 (94%)	13 (6%)	18	36
All	All	1153/1290 (89%)	1134 (98%)	19 (2%)	55	78

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ILE
1	A	191	GLU
1	B	21	LEU
1	D	191	GLU
1	D	255	LEU
1	D	258	GLU
1	E	28	LEU

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Mol	Chain	Res	Type
1	E	41	SER
1	E	43	GLN
1	E	112	ARG
1	E	133	ASP
1	E	204	THR
1	E	207	VAL
1	E	213	LEU
1	E	214	ILE
1	E	217	ARG
1	E	285	ASP
1	E	287	THR
1	E	292	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	268	ASN
1	C	235	HIS
1	D	43	GLN
1	E	31	ASN
1	E	99	GLN
1	E	158	ASN
1	E	194	HIS

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 15 ligands modelled in this entry, 15 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	274/296 (92%)	1.11	51 (18%) 3 2	43, 60, 88, 112	0
1	B	272/296 (91%)	1.23	53 (19%) 3 2	42, 66, 97, 127	0
1	C	271/296 (91%)	1.33	68 (25%) 1 1	44, 63, 103, 123	0
1	D	268/296 (90%)	2.02	101 (37%) 1 0	44, 69, 101, 115	0
1	E	270/296 (91%)	1.27	56 (20%) 2 2	44, 62, 94, 107	0
All	All	1355/1480 (91%)	1.39	329 (24%) 2 1	42, 64, 100, 127	0

All (329) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	145	GLU	14.9
1	E	290	HIS	11.2
1	B	49	GLY	9.7
1	C	48	LEU	9.5
1	D	106	ALA	9.2
1	A	16	TRP	9.0
1	C	47	GLN	8.7
1	D	144	GLU	8.2
1	D	146	ARG	8.0
1	B	210	ALA	8.0
1	E	206	PRO	7.7
1	C	20	VAL	7.6
1	B	23	LYS	7.4
1	B	211	TYR	7.4
1	E	204	THR	7.2
1	E	285	ASP	7.2
1	E	209	PHE	7.0
1	D	149	GLU	6.9
1	D	170	GLN	6.9
1	D	131	LYS	6.8

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Mol	Chain	Res	Type	RSRZ
1	B	209	PHE	6.7
1	D	233	ASP	6.6
1	D	234	LEU	6.6
1	B	47	GLN	6.5
1	B	22	SER	6.5
1	C	44	TRP	6.4
1	D	47	GLN	6.4
1	E	207	VAL	6.4
1	C	49	GLY	6.4
1	C	290	HIS	6.4
1	E	24	ILE	6.3
1	E	265	THR	6.3
1	E	211	TYR	6.1
1	C	22	SER	6.1
1	C	106	ALA	6.0
1	D	107	GLU	6.0
1	A	47	GLN	6.0
1	B	21	LEU	6.0
1	A	46	GLU	6.0
1	C	45	TYR	5.9
1	C	43	GLN	5.8
1	D	235	HIS	5.7
1	D	206	PRO	5.7
1	B	51	HIS	5.7
1	D	289	GLN	5.7
1	C	211	TYR	5.6
1	D	46	GLU	5.5
1	A	22	SER	5.5
1	D	42	TYR	5.5
1	E	213	LEU	5.4
1	D	77	SER	5.3
1	A	190	ASP	5.3
1	E	291	PRO	5.3
1	D	142	LEU	5.3
1	D	32	VAL	5.3
1	A	44	TRP	5.3
1	E	210	ALA	5.3
1	D	207	VAL	5.2
1	A	189	LEU	5.2
1	D	37	ILE	5.2
1	B	50	ILE	5.1
1	D	109	ASP	5.1

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Mol	Chain	Res	Type	RSRZ
1	C	268	ASN	5.1
1	D	22	SER	5.0
1	B	48	LEU	5.0
1	A	51	HIS	4.9
1	D	211	TYR	4.9
1	D	236	TYR	4.9
1	B	206	PRO	4.8
1	B	44	TRP	4.8
1	B	26	PHE	4.7
1	C	23	LYS	4.7
1	D	253	ASP	4.6
1	C	207	VAL	4.6
1	D	251	SER	4.5
1	D	33	LEU	4.5
1	D	44	TRP	4.5
1	E	264	GLY	4.5
1	A	203	ALA	4.5
1	D	173	GLU	4.4
1	D	143	PRO	4.4
1	B	207	VAL	4.4
1	E	44	TRP	4.4
1	B	292	LEU	4.4
1	D	176	LYS	4.4
1	D	166	ASN	4.4
1	D	260	GLU	4.4
1	E	33	LEU	4.4
1	D	108	HIS	4.3
1	E	36	ILE	4.3
1	C	184	LEU	4.3
1	E	41	SER	4.3
1	C	109	ASP	4.3
1	D	219	VAL	4.3
1	A	209	PHE	4.3
1	D	209	PHE	4.2
1	A	204	THR	4.2
1	C	209	PHE	4.2
1	D	36	ILE	4.2
1	D	265	THR	4.2
1	A	17	HIS	4.2
1	A	205	THR	4.2
1	D	263	PHE	4.2
1	D	23	LYS	4.2

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Mol	Chain	Res	Type	RSRZ
1	C	258	GLU	4.1
1	D	49	GLY	4.1
1	C	208	PRO	4.1
1	D	58	SER	4.1
1	E	212	THR	4.1
1	C	236	TYR	4.1
1	E	266	ALA	4.1
1	D	261	ASP	4.0
1	A	48	LEU	4.0
1	B	24	ILE	4.0
1	D	257	GLU	3.9
1	C	51	HIS	3.9
1	D	268	ASN	3.9
1	E	26	PHE	3.9
1	D	105	PRO	3.9
1	D	232	GLY	3.8
1	C	213	LEU	3.8
1	D	267	ALA	3.8
1	A	202	LEU	3.8
1	E	32	VAL	3.8
1	D	57	PHE	3.8
1	B	29	LEU	3.7
1	C	21	LEU	3.7
1	C	46	GLU	3.7
1	E	42	TYR	3.7
1	D	30	LEU	3.7
1	D	250	LEU	3.7
1	C	257	GLU	3.6
1	E	205	THR	3.6
1	E	131	LYS	3.6
1	C	37	ILE	3.6
1	B	46	GLU	3.6
1	B	58	SER	3.6
1	B	52	LEU	3.5
1	E	208	PRO	3.5
1	B	27	ARG	3.5
1	B	45	TYR	3.5
1	C	29	LEU	3.5
1	E	268	ASN	3.5
1	D	24	ILE	3.5
1	D	174	ALA	3.4
1	D	139	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	D	205	THR	3.4
1	D	215	LEU	3.4
1	D	266	ALA	3.4
1	C	42	TYR	3.4
1	C	41	SER	3.4
1	A	186	ASP	3.4
1	D	112	ARG	3.4
1	A	26	PHE	3.3
1	A	191	GLU	3.3
1	A	261	ASP	3.3
1	D	48	LEU	3.3
1	E	29	LEU	3.3
1	A	212	THR	3.3
1	E	235	HIS	3.3
1	A	268	ASN	3.2
1	B	56	PRO	3.2
1	C	79	PHE	3.2
1	D	51	HIS	3.2
1	D	148	THR	3.2
1	C	180	ILE	3.2
1	D	41	SER	3.2
1	D	252	ALA	3.1
1	B	220	TYR	3.1
1	D	288	GLY	3.1
1	C	265	THR	3.1
1	C	203	ALA	3.1
1	E	190	ASP	3.1
1	D	104	LEU	3.1
1	E	54	VAL	3.1
1	B	43	GLN	3.1
1	E	258	GLU	3.1
1	D	247	TYR	3.1
1	C	210	ALA	3.1
1	C	260	GLU	3.1
1	B	208	PRO	3.1
1	D	101	ARG	3.0
1	C	50	ILE	3.0
1	E	214	ILE	3.0
1	A	53	THR	3.0
1	D	129	LEU	3.0
1	C	289	GLN	3.0
1	C	267	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	29	LEU	2.9
1	D	255	LEU	2.9
1	B	205	THR	2.9
1	A	45	TYR	2.9
1	D	220	TYR	2.9
1	B	236	TYR	2.9
1	C	187	ASN	2.8
1	B	287	THR	2.8
1	A	23	LYS	2.8
1	B	214	ILE	2.8
1	D	191	GLU	2.8
1	E	289	GLN	2.8
1	A	40	ILE	2.8
1	D	45	TYR	2.8
1	D	50	ILE	2.8
1	A	32	VAL	2.8
1	B	59	LEU	2.8
1	A	50	ILE	2.7
1	A	213	LEU	2.7
1	A	200	GLU	2.7
1	C	103	ILE	2.7
1	E	168	ILE	2.7
1	B	83	ARG	2.7
1	B	111	HIS	2.7
1	D	210	ALA	2.7
1	E	48	LEU	2.7
1	B	148	THR	2.7
1	A	194	HIS	2.7
1	D	40	ILE	2.6
1	E	25	ILE	2.6
1	A	83	ARG	2.6
1	D	71	ARG	2.6
1	A	269	ASP	2.6
1	A	217	ARG	2.6
1	B	77	SER	2.6
1	C	204	THR	2.6
1	C	26	PHE	2.6
1	E	240	PHE	2.6
1	C	174	ALA	2.6
1	C	206	PRO	2.6
1	B	36	ILE	2.6
1	D	27	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	205	THR	2.5
1	C	220	TYR	2.5
1	A	197	GLY	2.5
1	E	232	GLY	2.5
1	C	25	ILE	2.5
1	B	60	LEU	2.5
1	D	212	THR	2.5
1	A	42	TYR	2.5
1	A	25	ILE	2.5
1	C	191	GLU	2.5
1	E	133	ASP	2.5
1	D	25	ILE	2.5
1	B	228	PHE	2.5
1	B	240	PHE	2.5
1	D	216	GLN	2.5
1	C	255	LEU	2.5
1	D	52	LEU	2.5
1	A	235	HIS	2.5
1	D	80	VAL	2.5
1	E	267	ALA	2.5
1	C	40	ILE	2.4
1	E	279	ILE	2.4
1	A	240	PHE	2.4
1	B	135	THR	2.4
1	C	108	HIS	2.4
1	D	264	GLY	2.4
1	A	43	GLN	2.4
1	B	145	GLU	2.4
1	D	81	GLU	2.4
1	D	39	ILE	2.4
1	E	40	ILE	2.4
1	B	217	ARG	2.4
1	C	202	LEU	2.4
1	C	266	ALA	2.4
1	C	107	GLU	2.4
1	E	114	ILE	2.3
1	E	261	ASP	2.3
1	A	267	ALA	2.3
1	C	77	SER	2.3
1	C	235	HIS	2.3
1	A	199	CYS	2.3
1	A	174	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	256	ALA	2.3
1	E	224	THR	2.3
1	B	151	LEU	2.3
1	D	180	ILE	2.3
1	E	255	LEU	2.3
1	B	42	TYR	2.3
1	C	281	ARG	2.3
1	D	182	TYR	2.3
1	E	37	ILE	2.2
1	E	39	ILE	2.2
1	E	275	MET	2.2
1	D	78	ARG	2.2
1	A	187	ASN	2.2
1	D	224	THR	2.2
1	A	255	LEU	2.2
1	A	270	LEU	2.2
1	C	24	ILE	2.2
1	D	179	ASP	2.2
1	B	32	VAL	2.2
1	A	277	ASN	2.2
1	A	247	TYR	2.2
1	B	238	THR	2.2
1	E	217	ARG	2.2
1	D	203	ALA	2.2
1	B	224	THR	2.2
1	E	278	THR	2.2
1	D	258	GLU	2.2
1	C	27	ARG	2.2
1	C	240	PHE	2.2
1	B	33	LEU	2.1
1	B	53	THR	2.1
1	D	213	LEU	2.1
1	D	68	LEU	2.1
1	E	189	LEU	2.1
1	B	37	ILE	2.1
1	D	228	PHE	2.1
1	E	231	VAL	2.1
1	A	252	ALA	2.1
1	C	256	ALA	2.1
1	B	40	ILE	2.1
1	D	208	PRO	2.1
1	E	194	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	263	PHE	2.1
1	D	110	ALA	2.1
1	C	101	ARG	2.1
1	A	131	LYS	2.1
1	A	77	SER	2.1
1	C	84	ASN	2.0
1	C	200	GLU	2.0
1	C	157	THR	2.0
1	A	19	SER	2.0
1	D	28	LEU	2.0
1	C	39	ILE	2.0
1	C	212	THR	2.0
1	C	190	ASP	2.0
1	B	235	HIS	2.0
1	E	220	TYR	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](#)

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	D	304	1/1	0.02	0.45	97,97,97,97	0
2	ZN	E	303	1/1	0.32	0.46	49,49,49,49	0
2	ZN	B	302	1/1	0.38	0.29	61,61,61,61	0
2	ZN	D	301	1/1	0.42	0.24	58,58,58,58	0
2	ZN	E	302	1/1	0.48	0.22	65,65,65,65	0
2	ZN	A	303	1/1	0.53	0.28	76,76,76,76	0
2	ZN	C	301	1/1	0.55	0.26	49,49,49,49	0
2	ZN	C	303	1/1	0.55	0.32	67,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	E	301	1/1	0.63	0.29	54,54,54,54	0
2	ZN	D	303	1/1	0.68	0.18	78,78,78,78	0
2	ZN	A	302	1/1	0.76	0.15	61,61,61,61	0
2	ZN	A	301	1/1	0.79	0.13	64,64,64,64	0
2	ZN	D	302	1/1	0.81	0.14	80,80,80,80	0
2	ZN	C	302	1/1	0.81	0.15	65,65,65,65	0
2	ZN	B	301	1/1	0.85	0.12	100,100,100,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.