Full wwPDB X-ray Structure Validation Report (i)

Jun 2, 2022 – 06:05 PM JST

PDB ID	:	6IV3
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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

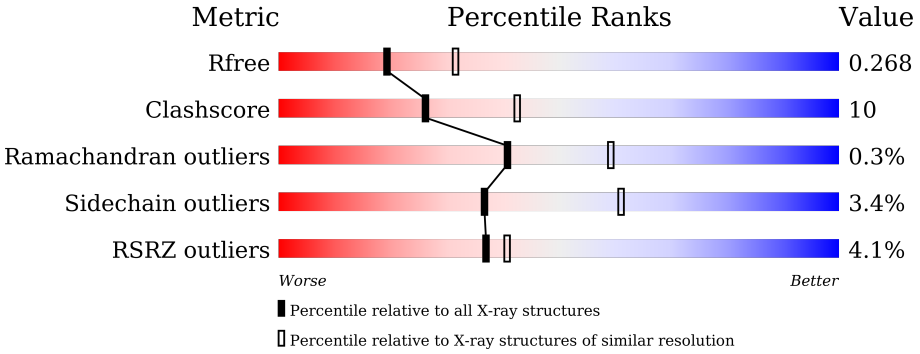
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.28.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.28.1

1 Overall quality at a glance (i)

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}	130704	5743 (2.54-2.50)
Clashscore	141614	6463 (2.54-2.50)
Ramachandran outliers	138981	6335 (2.54-2.50)
Sidechain outliers	138945	6337 (2.54-2.50)
RSRZ outliers	127900	5630 (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 <=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	2%, 71%, 21%, 7%
1	B	296	4%, 75%, 16%, • , 8%
1	C	296	5%, 74%, 18%, 8%
1	D	296	4%, 70%, 20%, • , 9%
1	E	296	3%, 69%, 19%, • , 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 (i)

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

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 (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is water.

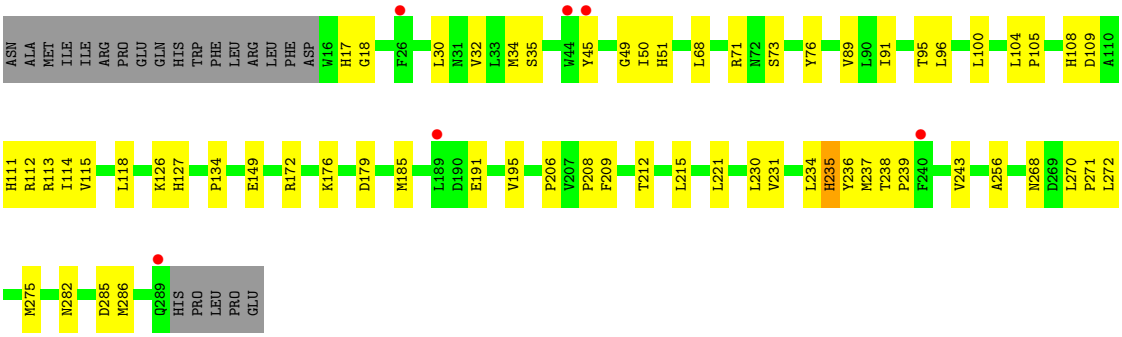
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	26	26	26	0	0
			Total	O		
3	B	13	13	13	0	0
			Total	O		
3	C	18	18	18	0	0
			Total	O		
3	D	8	8	8	0	0
			Total	O		
3	E	15	15	15	0	0

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.

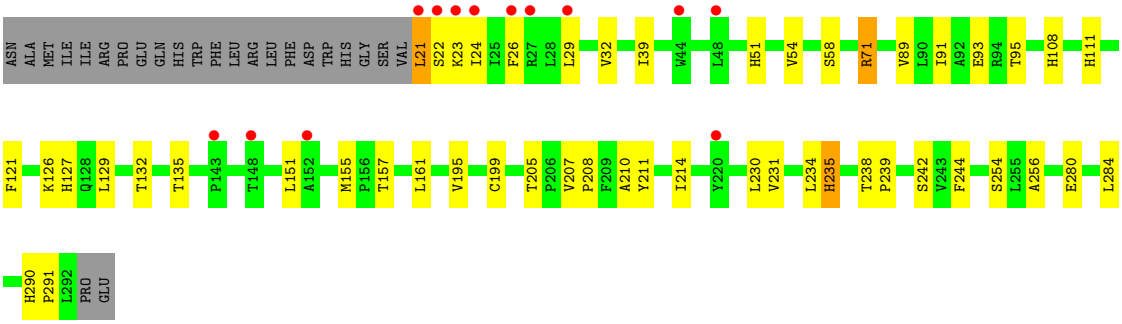
• Molecule 1: Bestrophin homolog

Chain A:



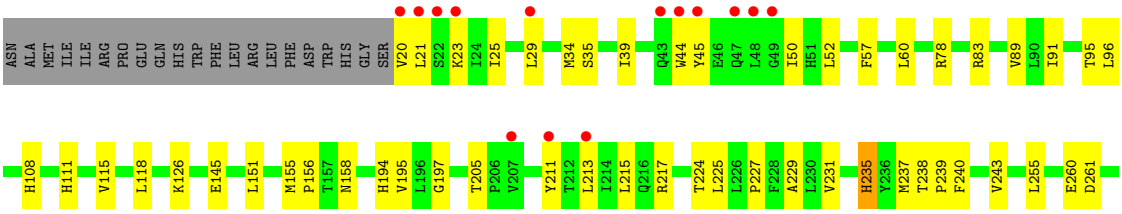
• Molecule 1: Bestrophin homolog

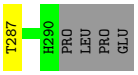
Chain B:



• Molecule 1: Bestrophin homolog

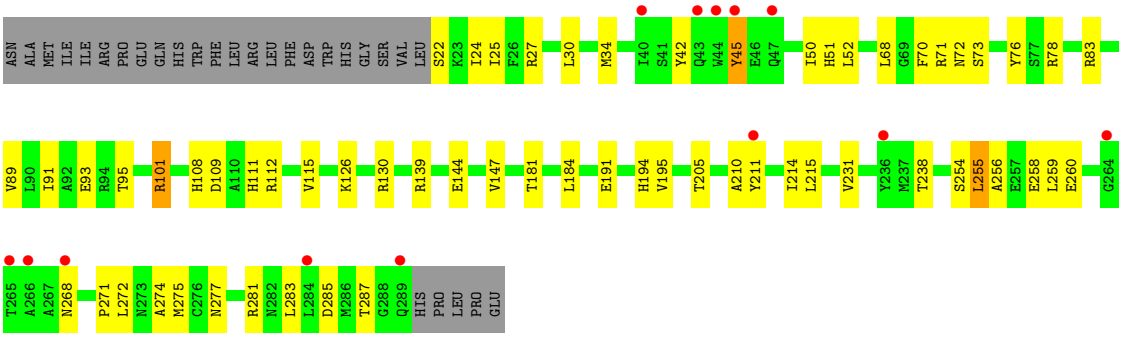
Chain C:





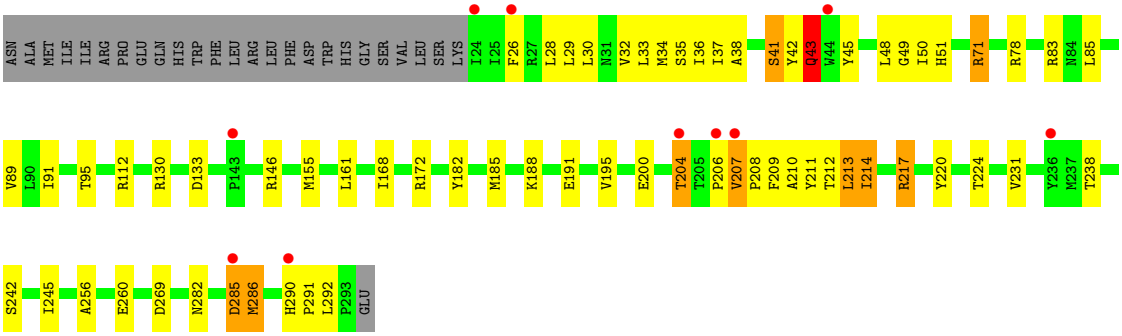
• Molecule 1: Bestrophin homolog

Chain D:



• Molecule 1: Bestrophin homolog

Chain E:



4 Data and refinement statistics (i)

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 – 2.51	Depositor EDS
% Data completeness (in resolution range)	98.7 (66.11-2.52) 98.6 (66.11-2.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.46 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.234 , 0.265 0.241 , 0.268	Depositor DCC
R_{free} test set	2006 reflections (2.02%)	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)	0.30 , 50.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o,F_c correlation	0.91	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.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: 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	0.49	0/2207	0.61	0/3006
1	B	0.43	0/2205	0.59	0/3001
1	C	0.43	0/2153	0.58	0/2932
1	D	0.45	0/2152	0.59	0/2930
1	E	0.49	0/2161	0.62	0/2945
All	All	0.46	0/10878	0.60	0/14814

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2160	0	2191	57	0
1	B	2159	0	2212	39	0
1	C	2112	0	2152	41	0
1	D	2109	0	2143	53	0
1	E	2117	0	2155	65	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
2	C	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	1	0
3	D	8	0	0	1	0
3	E	15	0	0	2	0
All	All	10752	0	10853	223	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 10.

All (223) 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:E:168:ILE:HD12	1:E:185:MET:HE3	1.60	0.83
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
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: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:B:205:THR:HG21	1:D:83:ARG:HE	1.56	0.70
1:A:268:ASN:ND2	1:E:210:ALA:CB	2.54	0.69
1:C:21:LEU:C	1:C:21:LEU:HD13	2.13	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:91:ILE:O	1:A:95:THR:HG23	1.97	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: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:C:205:THR:OG1	1:E:83:ARG:NH1	2.35	0.59
1:C:20:VAL:HG13	1:C:23:LYS:H	1.69	0.58
1:E:200:GLU:O	1:E:204:THR:OG1	2.20	0.58
1:A:109:ASP:OD1	1:A:112:ARG:NH2	2.37	0.58
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: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: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:207:VAL:O	1:E:208:PRO:C	2.43	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:D:271:PRO:HD2	1:D:275:MET:HE3	1.88	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:E:214:ILE:C	1:E:214:ILE:HD13	2.28	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:57:PHE:HE2	1:C:225:LEU:HB3	1.73	0.54
1:A:271:PRO:HD2	1:A:275:MET:HE2	1.90	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:270:LEU:HB3	1:A:275:MET:HE1	1.91	0.52
1:A:272:LEU:HD23	1:A:275:MET:HE1	1.90	0.52
1:A:285:ASP:HB3	1:E:155: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:A:270:LEU:HB3	1:A:275:MET:CE	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:E:290:HIS:N	1:E:291:PRO:CD	2.73	0.51
1:C:235:HIS:CD2	1:D:51:HIS:CE1	2.98	0.51
1:D:255:LEU:O	1:D:255:LEU:HG	2.11	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:D:108:HIS:HD1	1:D:111:HIS:CE1	2.30	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:C:145:GLU:H	1:C:145:GLU:CD	2.15	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:B:127:HIS:ND1	1:B:132:THR:OG1	2.38	0.49
1:A:32:VAL:HG23	1:A:243:VAL:HG21	1.94	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:B:155:MET:HE1	1:D:285:ASP:HB3	1.95	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:101:ARG:NH1	1:D:108:HIS:CE1	2.81	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:B:135:THR:HG22	1:B:151:LEU:CD1	2.45	0.47
1:C:21:LEU:C	1:C:21:LEU:CD1	2.83	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: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: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:237:MET:HE2	1:D:52:LEU:HD21	1.98	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:E:211:TYR:C	1:E:214:ILE:CG2	2.86	0.43
1:A:104:LEU:HD21	1:A:176:LYS:O	2.18	0.43
1:A:272:LEU:HD23	1:A:275:MET:CE	2.49	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:A:114:ILE:O	1:A:118:LEU:HG	2.18	0.43
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: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:E:220:TYR:O	1:E:224:THR:HG23	2.18	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: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:A:100:LEU:HG	1:A:185:MET:CE	2.50	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:237:MET:HE3	1:A:237:MET:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:96:LEU:HD23	1:C:118:LEU:HD21	2.02	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:B:54:VAL:HG22	1:B:230:LEU:HD23	2.02	0.41
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:236:TYR:HB2	1:E:50:ILE:HD13	2.03	0.41
1:A:285:ASP:CB	1:E:155:MET:HE1	2.51	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:A:282:ASN:HA	1:E:155:MET:HE1	2.03	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%)	22	37
1	B	270/296 (91%)	262 (97%)	7 (3%)	1 (0%)	34	53
1	C	269/296 (91%)	265 (98%)	4 (2%)	0	100	100
1	D	266/296 (90%)	257 (97%)	9 (3%)	0	100	100
1	E	268/296 (90%)	264 (98%)	3 (1%)	1 (0%)	34	53
All	All	1345/1480 (91%)	1312 (98%)	29 (2%)	4 (0%)	41	59

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 (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	233/258 (90%)	227 (97%)	6 (3%)	46	70
1	B	236/258 (92%)	229 (97%)	7 (3%)	41	66
1	C	227/258 (88%)	225 (99%)	2 (1%)	78	91
1	D	228/258 (88%)	220 (96%)	8 (4%)	36	60

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	229/258 (89%)	213 (93%)	16 (7%)	15	27
All	All	1153/1290 (89%)	1114 (97%)	39 (3%)	37	61

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	HIS
1	A	51	HIS
1	A	73	SER
1	A	126	LYS
1	A	191	GLU
1	A	235	HIS
1	B	21	LEU
1	B	51	HIS
1	B	58	SER
1	B	71	ARG
1	B	126	LYS
1	B	211	TYR
1	B	254	SER
1	C	126	LYS
1	C	235	HIS
1	D	45	TYR
1	D	101	ARG
1	D	126	LYS
1	D	144	GLU
1	D	191	GLU
1	D	255	LEU
1	D	258	GLU
1	D	268	ASN
1	E	26	PHE
1	E	28	LEU
1	E	41	SER
1	E	43	GLN
1	E	71	ARG
1	E	112	ARG
1	E	133	ASP
1	E	191	GLU
1	E	204	THR
1	E	207	VAL
1	E	213	LEU
1	E	214	ILE
1	E	217	ARG

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Mol	Chain	Res	Type
1	E	285	ASP
1	E	286	MET
1	E	292	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	HIS
1	A	268	ASN
1	C	235	HIS
1	E	194	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

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 (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	274/296 (92%)	-0.03	6 (2%)	62	65	43, 60, 88, 112	0
1	B	272/296 (91%)	0.14	13 (4%)	30	33	42, 66, 97, 127	0
1	C	271/296 (91%)	0.16	14 (5%)	27	29	44, 63, 103, 123	0
1	D	268/296 (90%)	0.25	13 (4%)	29	32	44, 69, 101, 115	0
1	E	270/296 (91%)	0.08	10 (3%)	41	45	44, 62, 94, 107	0
All	All	1355/1480 (91%)	0.12	56 (4%)	37	41	42, 64, 100, 127	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	289	GLN	6.5
1	C	48	LEU	6.2
1	C	44	TRP	6.0
1	B	23	LYS	5.8
1	B	26	PHE	5.6
1	B	24	ILE	5.5
1	D	236	TYR	5.3
1	C	45	TYR	5.0
1	D	266	ALA	4.9
1	C	47	GLN	4.6
1	B	22	SER	4.5
1	D	265	THR	4.4
1	D	44	TRP	4.3
1	B	21	LEU	4.2
1	E	26	PHE	4.2
1	C	211	TYR	4.0
1	B	220	TYR	4.0
1	C	49	GLY	4.0
1	B	143	PRO	3.7
1	C	22	SER	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	44	TRP	3.7
1	C	23	LYS	3.6
1	E	290	HIS	3.6
1	D	40	ILE	3.4
1	E	285	ASP	3.3
1	E	207	VAL	3.2
1	B	27	ARG	3.2
1	E	24	ILE	3.0
1	E	44	TRP	3.0
1	A	189	LEU	2.9
1	B	48	LEU	2.8
1	C	20	VAL	2.7
1	E	204	THR	2.6
1	B	152	ALA	2.6
1	A	45	TYR	2.5
1	D	268	ASN	2.5
1	B	44	TRP	2.5
1	D	45	TYR	2.5
1	D	264	GLY	2.5
1	C	21	LEU	2.4
1	D	47	GLN	2.4
1	E	143	PRO	2.4
1	C	29	LEU	2.4
1	D	43	GLN	2.3
1	D	211	TYR	2.3
1	A	289	GLN	2.3
1	C	43	GLN	2.3
1	A	240	PHE	2.2
1	C	207	VAL	2.2
1	D	284	LEU	2.2
1	B	148	THR	2.1
1	C	213	LEU	2.1
1	B	29	LEU	2.1
1	A	26	PHE	2.0
1	E	236	TYR	2.0
1	E	206	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 monosaccharides 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	B	302	1/1	-0.24	0.24	61,61,61,61	0
2	ZN	E	302	1/1	-0.23	0.23	65,65,65,65	0
2	ZN	D	304	1/1	-0.08	0.49	97,97,97,97	0
2	ZN	C	301	1/1	0.37	0.22	49,49,49,49	0
2	ZN	C	303	1/1	0.46	0.27	67,67,67,67	0
2	ZN	D	303	1/1	0.48	0.16	78,78,78,78	0
2	ZN	A	303	1/1	0.56	0.28	76,76,76,76	0
2	ZN	B	301	1/1	0.56	0.09	100,100,100,100	0
2	ZN	C	302	1/1	0.73	0.11	65,65,65,65	0
2	ZN	E	301	1/1	0.74	0.24	54,54,54,54	0
2	ZN	E	303	1/1	0.77	0.60	49,49,49,49	0
2	ZN	D	301	1/1	0.80	0.31	58,58,58,58	0
2	ZN	A	301	1/1	0.80	0.10	64,64,64,64	0
2	ZN	D	302	1/1	0.88	0.14	80,80,80,80	0
2	ZN	A	302	1/1	0.89	0.12	61,61,61,61	0

6.5 Other polymers (i)

There are no such residues in this entry.