



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:15 PM GMT

PDB ID : 4YEU
Title : ELIC-GLIC chimera in the resting conformation
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Deposited on : 2015-02-24
Resolution : 4.60 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20026982

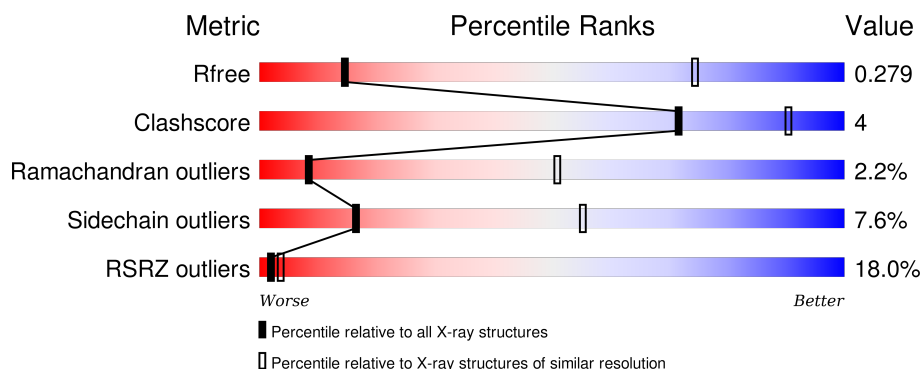
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.60 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}	91344	1089 (5.52-3.60)
Clashscore	102246	1004 (5.52-3.64)
Ramachandran outliers	100387	1131 (5.52-3.60)
Sidechain outliers	100360	1112 (5.50-3.60)
RSRZ outliers	91569	1092 (5.52-3.60)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	312	<div> <div>13%</div> <div>81%</div> <div>17%</div> <div>.</div> </div>
1	B	312	<div> <div>18%</div> <div>80%</div> <div>19%</div> <div>.</div> </div>
1	C	312	<div> <div>21%</div> <div>80%</div> <div>18%</div> <div>.</div> </div>
1	D	312	<div> <div>26%</div> <div>82%</div> <div>16%</div> <div>.</div> </div>
1	E	312	<div> <div>12%</div> <div>81%</div> <div>17%</div> <div>.</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 12825 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 Cys-loop ligand-gated ion channel,Proton-gated ion channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	312	Total	C	N	O	S	0	0	0
			2565	1686	417	457	5			
1	B	312	Total	C	N	O	S	0	0	0
			2565	1686	417	457	5			
1	C	312	Total	C	N	O	S	0	0	0
			2565	1686	417	457	5			
1	D	312	Total	C	N	O	S	0	0	0
			2565	1686	417	457	5			
1	E	312	Total	C	N	O	S	0	0	0
			2565	1686	417	457	5			

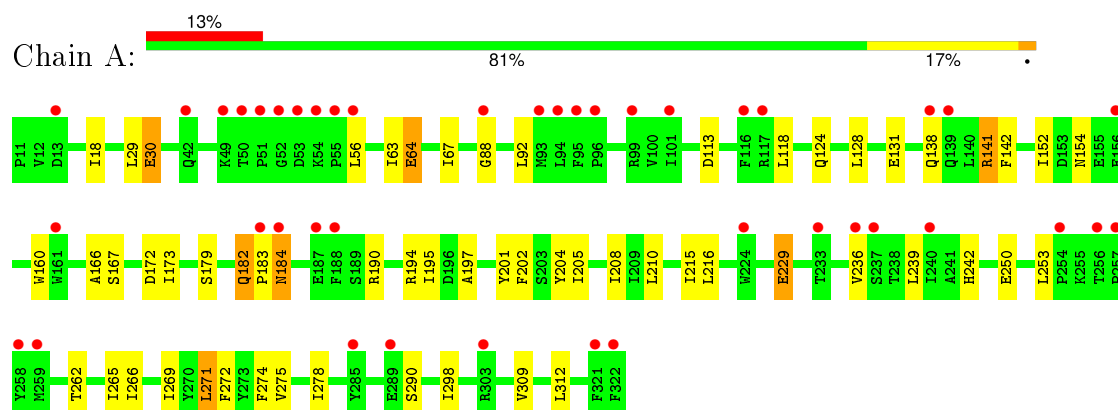
There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	164	GLY	-	insertion	UNP P0C7B7
B	164	GLY	-	insertion	UNP P0C7B7
C	164	GLY	-	insertion	UNP P0C7B7
D	164	GLY	-	insertion	UNP P0C7B7
E	164	GLY	-	insertion	UNP P0C7B7

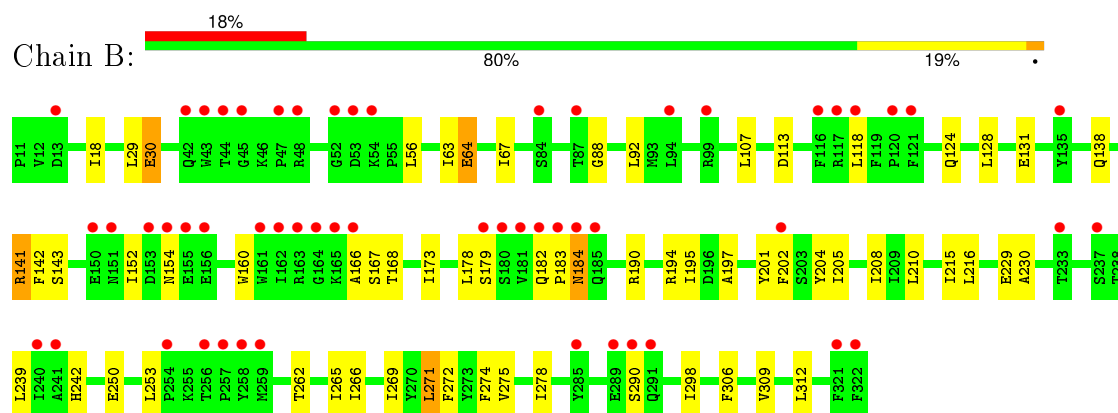
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

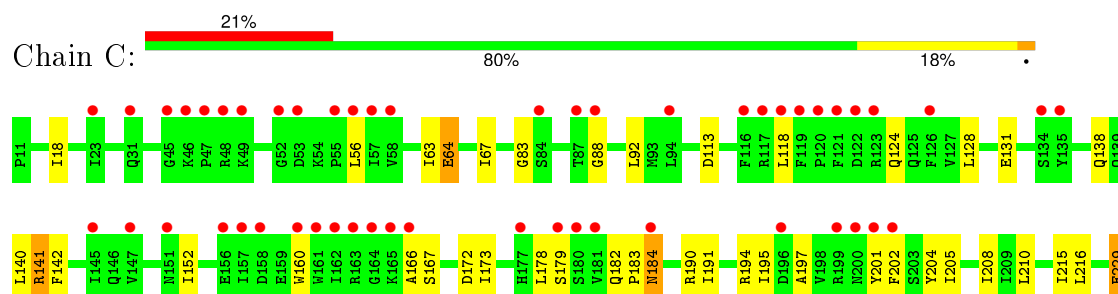
- Molecule 1: Cys-loop ligand-gated ion channel,Proton-gated ion channel

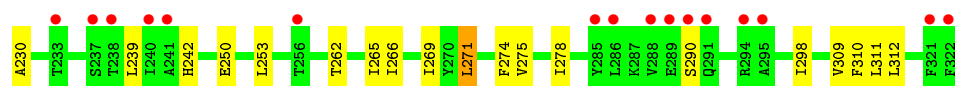


- Molecule 1: Cys-loop ligand-gated ion channel,Proton-gated ion channel

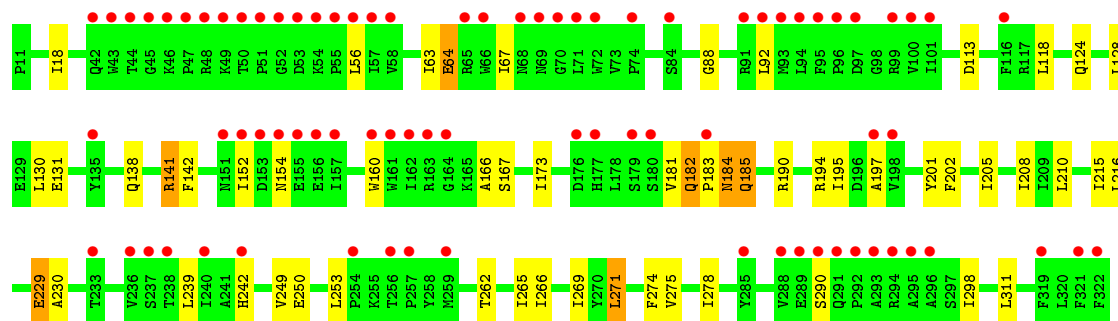
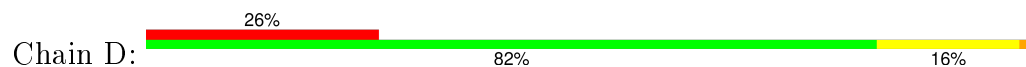


- Molecule 1: Cys-loop ligand-gated ion channel,Proton-gated ion channel

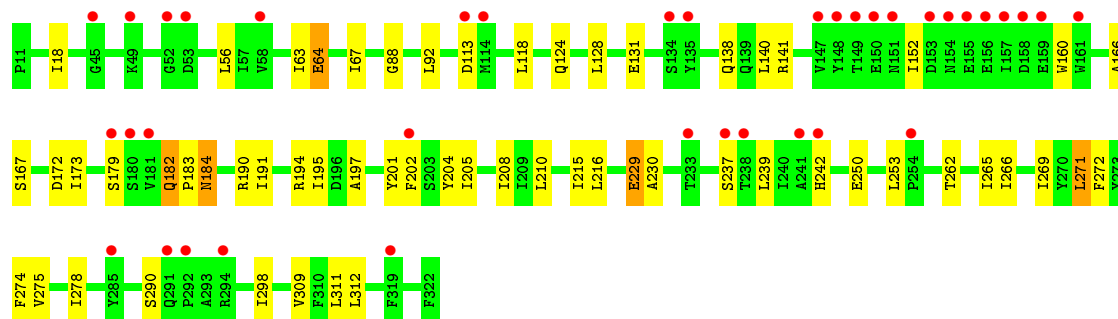
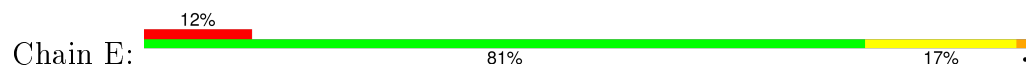




- Molecule 1: Cys-loop ligand-gated ion channel,Proton-gated ion channel



- Molecule 1: Cys-loop ligand-gated ion channel,Proton-gated ion channel



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	132.44Å 217.96Å 229.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 4.60 49.23 – 4.60	Depositor EDS
% Data completeness (in resolution range)	99.3 (20.00-4.60) 99.3 (49.23-4.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 4.64Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.242 , 0.248 0.277 , 0.279	Depositor DCC
R_{free} test set	946 reflections (5.13%)	DCC
Wilson B-factor (Å ²)	254.6	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 102.1	EDS
Estimated twinning fraction	0.016 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.024 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	3 of 18685 reflections (0.016%)	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	12825	wwPDB-VP
Average B, all atoms (Å ²)	177.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.38	0/2638	0.61	0/3600
1	B	0.38	0/2638	0.61	0/3600
1	C	0.38	0/2638	0.61	0/3600
1	D	0.38	0/2638	0.63	0/3600
1	E	0.38	0/2638	0.61	0/3600
All	All	0.38	0/13190	0.62	0/18000

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	2565	0	2545	23	0
1	B	2565	0	2545	26	0
1	C	2565	0	2545	24	0
1	D	2565	0	2545	24	0
1	E	2565	0	2545	23	0
All	All	12825	0	12725	109	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (109) 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:208:ILE:HD12	1:D:249:VAL:HG22	1.46	0.97
1:D:182:GLN:HB3	1:D:185:GLN:HB3	1.57	0.85
1:D:167:SER:HB3	1:D:194:ARG:HG3	1.71	0.73
1:E:167:SER:HB3	1:E:194:ARG:HG3	1.71	0.72
1:C:167:SER:HB3	1:C:194:ARG:HG3	1.71	0.72
1:B:167:SER:HB3	1:B:194:ARG:HG3	1.72	0.72
1:A:167:SER:HB3	1:A:194:ARG:HG3	1.71	0.70
1:D:182:GLN:HG3	1:D:184:ASN:H	1.59	0.67
1:C:215:ILE:HD12	1:C:242:HIS:HA	1.82	0.61
1:D:215:ILE:HD12	1:D:242:HIS:HA	1.82	0.61
1:A:215:ILE:HD12	1:A:242:HIS:HA	1.81	0.61
1:B:215:ILE:HD12	1:B:242:HIS:HA	1.83	0.60
1:E:215:ILE:HD12	1:E:242:HIS:HA	1.83	0.60
1:A:166:ALA:HB2	1:A:195:ILE:HG12	1.85	0.58
1:C:166:ALA:HB2	1:C:195:ILE:HG12	1.87	0.57
1:A:239:LEU:HD21	1:E:215:ILE:HG22	1.87	0.56
1:B:166:ALA:HB2	1:B:195:ILE:HG12	1.86	0.56
1:E:166:ALA:HB2	1:E:195:ILE:HG12	1.86	0.56
1:D:166:ALA:HB2	1:D:195:ILE:HG12	1.87	0.55
1:B:215:ILE:HG22	1:C:239:LEU:HD21	1.89	0.54
1:C:215:ILE:HG22	1:D:239:LEU:HD21	1.88	0.54
1:A:229:GLU:HG3	1:E:230:ALA:HB2	1.88	0.54
1:A:274:PHE:O	1:A:278:ILE:HG12	2.08	0.53
1:A:215:ILE:HG22	1:B:239:LEU:HD21	1.89	0.53
1:E:274:PHE:O	1:E:278:ILE:HG12	2.09	0.53
1:B:274:PHE:O	1:B:278:ILE:HG12	2.08	0.53
1:C:274:PHE:O	1:C:278:ILE:HG12	2.09	0.53
1:D:274:PHE:O	1:D:278:ILE:HG12	2.10	0.51
1:C:216:LEU:HD13	1:C:269:ILE:HG23	1.93	0.50
1:B:216:LEU:HD13	1:B:269:ILE:HG23	1.94	0.49
1:B:178:LEU:HD22	1:B:182:GLN:NE2	2.28	0.49
1:D:216:LEU:HD13	1:D:269:ILE:HG23	1.94	0.49
1:A:216:LEU:HD13	1:A:269:ILE:HG23	1.94	0.49
1:D:215:ILE:HG22	1:E:239:LEU:HD21	1.95	0.48
1:E:265:ILE:O	1:E:269:ILE:HG13	2.13	0.48
1:E:216:LEU:HD13	1:E:269:ILE:HG23	1.94	0.48
1:A:266:ILE:HA	1:A:269:ILE:HD12	1.95	0.48
1:B:266:ILE:HA	1:B:269:ILE:HD12	1.96	0.48
1:D:266:ILE:HA	1:D:269:ILE:HD12	1.96	0.48
1:A:265:ILE:O	1:A:269:ILE:HG13	2.14	0.47
1:B:160:TRP:HB3	1:B:197:ALA:HB1	1.96	0.47
1:C:266:ILE:HA	1:C:269:ILE:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:266:ILE:HA	1:E:269:ILE:HD12	1.96	0.47
1:B:131:GLU:HB2	1:B:190:ARG:HG3	1.96	0.47
1:A:160:TRP:HB3	1:A:197:ALA:HB1	1.96	0.47
1:D:160:TRP:HB3	1:D:197:ALA:HB1	1.96	0.47
1:E:64:GLU:HA	1:E:67:ILE:HD12	1.97	0.47
1:E:131:GLU:HB2	1:E:190:ARG:HG3	1.97	0.47
1:D:265:ILE:O	1:D:269:ILE:HG13	2.14	0.47
1:C:230:ALA:HB2	1:D:229:GLU:HG3	1.97	0.47
1:E:160:TRP:HB3	1:E:197:ALA:HB1	1.97	0.47
1:A:131:GLU:HB2	1:A:190:ARG:HG3	1.96	0.47
1:D:131:GLU:HB2	1:D:190:ARG:HG3	1.96	0.47
1:C:131:GLU:HB2	1:C:190:ARG:HG3	1.97	0.46
1:A:64:GLU:HA	1:A:67:ILE:HD12	1.97	0.46
1:A:271:LEU:O	1:A:275:VAL:HG23	2.15	0.46
1:B:271:LEU:O	1:B:275:VAL:HG23	2.16	0.46
1:C:265:ILE:O	1:C:269:ILE:HG13	2.16	0.46
1:B:173:ILE:HD13	1:B:190:ARG:HB3	1.97	0.46
1:C:64:GLU:HA	1:C:67:ILE:HD12	1.96	0.46
1:B:265:ILE:O	1:B:269:ILE:HG13	2.14	0.46
1:C:173:ILE:HD13	1:C:190:ARG:HB3	1.97	0.46
1:E:271:LEU:O	1:E:275:VAL:HG23	2.16	0.46
1:D:64:GLU:HA	1:D:67:ILE:HD12	1.97	0.46
1:A:182:GLN:N	1:A:183:PRO:HD3	2.31	0.46
1:A:204:TYR:HD1	1:A:208:ILE:HD12	1.80	0.46
1:D:173:ILE:HD13	1:D:190:ARG:HB3	1.98	0.46
1:B:107:LEU:HD13	1:C:83:GLY:H	1.81	0.45
1:A:173:ILE:HD13	1:A:190:ARG:HB3	1.97	0.45
1:D:271:LEU:O	1:D:275:VAL:HG23	2.16	0.45
1:D:63:ILE:HG12	1:D:92:LEU:HD12	1.97	0.45
1:C:160:TRP:HB3	1:C:197:ALA:HB1	1.97	0.45
1:C:178:LEU:HD22	1:C:182:GLN:NE2	2.30	0.45
1:C:182:GLN:N	1:C:183:PRO:HD3	2.31	0.45
1:E:173:ILE:HD13	1:E:190:ARG:HB3	1.98	0.45
1:D:230:ALA:HB2	1:E:229:GLU:HG3	1.98	0.45
1:B:204:TYR:HD1	1:B:208:ILE:HD12	1.82	0.45
1:B:182:GLN:N	1:B:183:PRO:HD3	2.32	0.45
1:B:29:LEU:HD23	1:B:30:GLU:HG2	1.99	0.45
1:A:236:VAL:HG11	1:E:237:SER:HB2	1.99	0.45
1:C:271:LEU:O	1:C:275:VAL:HG23	2.18	0.44
1:B:230:ALA:HB2	1:C:229:GLU:HG3	1.99	0.44
1:A:63:ILE:HG12	1:A:92:LEU:HD12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:204:TYR:HD1	1:E:208:ILE:HD12	1.81	0.44
1:B:63:ILE:HG12	1:B:92:LEU:HD12	1.99	0.44
1:D:182:GLN:HA	1:D:183:PRO:HD3	1.96	0.44
1:E:182:GLN:N	1:E:183:PRO:HD3	2.32	0.44
1:B:64:GLU:HA	1:B:67:ILE:HD12	1.98	0.44
1:B:306:PHE:HA	1:B:309:VAL:HG12	2.00	0.43
1:C:204:TYR:HD1	1:C:208:ILE:HD12	1.82	0.43
1:E:63:ILE:HG12	1:E:92:LEU:HD12	1.99	0.43
1:A:29:LEU:HD23	1:A:30:GLU:HG2	2.00	0.43
1:C:63:ILE:HG12	1:C:92:LEU:HD12	1.99	0.42
1:C:141:ARG:HE	1:C:142:PHE:HE2	1.68	0.41
1:B:272:PHE:CE2	1:B:309:VAL:HG13	2.55	0.41
1:D:141:ARG:HE	1:D:142:PHE:HE2	1.68	0.41
1:D:182:GLN:CB	1:D:185:GLN:HB3	2.41	0.41
1:C:309:VAL:HA	1:C:312:LEU:HD12	2.03	0.41
1:B:143:SER:HB3	1:B:168:THR:HG23	2.03	0.41
1:E:309:VAL:HA	1:E:312:LEU:HD12	2.03	0.41
1:D:184:ASN:HD22	1:D:184:ASN:HA	1.58	0.41
1:E:140:LEU:HD13	1:E:191:ILE:HG13	2.03	0.41
1:B:309:VAL:HA	1:B:312:LEU:HD12	2.03	0.40
1:B:141:ARG:HE	1:B:142:PHE:HE2	1.67	0.40
1:A:272:PHE:HZ	1:A:309:VAL:HG12	1.85	0.40
1:A:309:VAL:HA	1:A:312:LEU:HD12	2.03	0.40
1:E:272:PHE:HZ	1:E:309:VAL:HG12	1.87	0.40
1:A:141:ARG:HE	1:A:142:PHE:HE2	1.68	0.40
1:C:140:LEU:HD13	1:C:191:ILE:HG13	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	310/312 (99%)	278 (90%)	25 (8%)	7 (2%)	8	50
1	B	310/312 (99%)	278 (90%)	25 (8%)	7 (2%)	8	50
1	C	310/312 (99%)	278 (90%)	25 (8%)	7 (2%)	8	50
1	D	310/312 (99%)	280 (90%)	24 (8%)	6 (2%)	10	53
1	E	310/312 (99%)	278 (90%)	25 (8%)	7 (2%)	8	50
All	All	1550/1560 (99%)	1392 (90%)	124 (8%)	34 (2%)	8	51

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	250	GLU
1	B	250	GLU
1	C	250	GLU
1	D	250	GLU
1	E	250	GLU
1	E	141	ARG
1	A	141	ARG
1	A	202	PHE
1	A	290	SER
1	B	141	ARG
1	B	179	SER
1	B	202	PHE
1	B	290	SER
1	C	141	ARG
1	C	202	PHE
1	D	141	ARG
1	D	202	PHE
1	D	290	SER
1	E	202	PHE
1	E	290	SER
1	A	184	ASN
1	B	184	ASN
1	C	179	SER
1	C	290	SER
1	A	179	SER
1	C	184	ASN
1	E	179	SER
1	E	184	ASN
1	B	88	GLY
1	C	88	GLY
1	D	88	GLY

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Mol	Chain	Res	Type
1	E	88	GLY
1	A	88	GLY
1	D	181	VAL

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	283/283 (100%)	261 (92%)	22 (8%)	16	54
1	B	283/283 (100%)	263 (93%)	20 (7%)	18	57
1	C	283/283 (100%)	262 (93%)	21 (7%)	17	56
1	D	283/283 (100%)	260 (92%)	23 (8%)	15	52
1	E	283/283 (100%)	262 (93%)	21 (7%)	17	56
All	All	1415/1415 (100%)	1308 (92%)	107 (8%)	16	55

All (107) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ILE
1	A	30	GLU
1	A	56	LEU
1	A	64	GLU
1	A	113	ASP
1	A	118	LEU
1	A	124	GLN
1	A	128	LEU
1	A	138	GLN
1	A	152	ILE
1	A	154	ASN
1	A	172	ASP
1	A	182	GLN
1	A	184	ASN
1	A	201	TYR
1	A	205	ILE

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Mol	Chain	Res	Type
1	A	210	LEU
1	A	229	GLU
1	A	253	LEU
1	A	262	THR
1	A	271	LEU
1	A	298	ILE
1	B	18	ILE
1	B	30	GLU
1	B	56	LEU
1	B	64	GLU
1	B	113	ASP
1	B	118	LEU
1	B	124	GLN
1	B	128	LEU
1	B	138	GLN
1	B	152	ILE
1	B	154	ASN
1	B	184	ASN
1	B	201	TYR
1	B	205	ILE
1	B	210	LEU
1	B	229	GLU
1	B	253	LEU
1	B	262	THR
1	B	271	LEU
1	B	298	ILE
1	C	18	ILE
1	C	56	LEU
1	C	64	GLU
1	C	113	ASP
1	C	118	LEU
1	C	124	GLN
1	C	128	LEU
1	C	138	GLN
1	C	152	ILE
1	C	172	ASP
1	C	184	ASN
1	C	201	TYR
1	C	205	ILE
1	C	210	LEU
1	C	229	GLU
1	C	253	LEU

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Mol	Chain	Res	Type
1	C	262	THR
1	C	271	LEU
1	C	298	ILE
1	C	310	PHE
1	C	311	LEU
1	D	18	ILE
1	D	56	LEU
1	D	64	GLU
1	D	113	ASP
1	D	118	LEU
1	D	124	GLN
1	D	128	LEU
1	D	130	LEU
1	D	138	GLN
1	D	152	ILE
1	D	154	ASN
1	D	182	GLN
1	D	184	ASN
1	D	185	GLN
1	D	201	TYR
1	D	205	ILE
1	D	210	LEU
1	D	229	GLU
1	D	253	LEU
1	D	262	THR
1	D	271	LEU
1	D	298	ILE
1	D	311	LEU
1	E	18	ILE
1	E	56	LEU
1	E	64	GLU
1	E	113	ASP
1	E	118	LEU
1	E	124	GLN
1	E	128	LEU
1	E	138	GLN
1	E	152	ILE
1	E	172	ASP
1	E	182	GLN
1	E	184	ASN
1	E	201	TYR
1	E	205	ILE

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Mol	Chain	Res	Type
1	E	210	LEU
1	E	229	GLU
1	E	253	LEU
1	E	262	THR
1	E	271	LEU
1	E	298	ILE
1	E	311	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	ASN
1	A	200	ASN
1	A	252	ASN
1	B	112	ASN
1	B	182	GLN
1	C	112	ASN
1	C	200	ASN
1	C	252	ASN
1	D	112	ASN
1	D	184	ASN
1	D	200	ASN
1	D	252	ASN
1	E	112	ASN
1	E	200	ASN
1	E	252	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	312/312 (100%)	0.63	42 (13%) 4 5	73, 153, 255, 278	0
1	B	312/312 (100%)	0.70	55 (17%) 2 4	93, 191, 292, 300	0
1	C	312/312 (100%)	0.90	67 (21%) 1 3	109, 206, 298, 300	0
1	D	312/312 (100%)	1.06	80 (25%) 1 2	95, 172, 264, 294	0
1	E	312/312 (100%)	0.59	37 (11%) 6 7	80, 142, 231, 278	0
All	All	1560/1560 (100%)	0.78	281 (18%) 2 3	73, 170, 274, 300	0

All (281) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	53	ASP	10.0
1	B	183	PRO	9.2
1	E	156	GLU	8.5
1	D	53	ASP	7.9
1	D	49	LYS	7.3
1	E	53	ASP	6.9
1	E	157	ILE	6.8
1	B	53	ASP	6.4
1	C	322	PHE	6.0
1	E	150	GLU	6.0
1	B	180	SER	5.9
1	D	45	GLY	5.8
1	D	48	ARG	5.6
1	D	289	GLU	5.5
1	B	117	ARG	5.3
1	E	149	THR	5.3
1	C	48	ARG	5.3
1	C	165	LYS	5.3
1	B	156	GLU	5.2
1	B	184	ASN	5.2

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Mol	Chain	Res	Type	RSRZ
1	D	66	TRP	5.2
1	A	184	ASN	5.2
1	E	180	SER	5.1
1	B	179	SER	5.1
1	D	56	LEU	5.0
1	D	163	ARG	5.0
1	C	53	ASP	4.9
1	C	121	PHE	4.9
1	D	254	PRO	4.8
1	E	179	SER	4.8
1	D	94	LEU	4.8
1	C	294	ARG	4.7
1	D	69	ASN	4.7
1	D	52	GLY	4.6
1	D	161	TRP	4.6
1	E	151	ASN	4.6
1	D	153	ASP	4.6
1	C	179	SER	4.5
1	C	164	GLY	4.5
1	D	58	VAL	4.4
1	A	259	MET	4.4
1	A	183	PRO	4.4
1	D	151	ASN	4.4
1	B	163	ARG	4.4
1	D	91	ARG	4.4
1	B	164	GLY	4.3
1	D	72	TRP	4.3
1	D	71	LEU	4.3
1	A	55	PRO	4.3
1	D	47	PRO	4.2
1	D	155	GLU	4.2
1	C	58	VAL	4.2
1	E	158	ASP	4.2
1	C	180	SER	4.2
1	E	237	SER	4.2
1	D	154	ASN	4.2
1	B	289	GLU	4.1
1	C	147	VAL	4.1
1	A	52	GLY	4.1
1	D	156	GLU	4.1
1	C	156	GLU	4.1
1	D	46	LYS	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	94	LEU	4.1
1	D	55	PRO	4.0
1	B	44	THR	4.0
1	D	285	TYR	4.0
1	C	94	LEU	4.0
1	C	321	PHE	4.0
1	D	294	ARG	3.9
1	A	322	PHE	3.9
1	C	201	TYR	3.9
1	C	120	PRO	3.9
1	B	153	ASP	3.9
1	E	202	PHE	3.9
1	A	187	GLU	3.8
1	C	199	ARG	3.8
1	C	162	ILE	3.8
1	D	57	ILE	3.8
1	C	289	GLU	3.8
1	A	156	GLU	3.7
1	C	160	TRP	3.7
1	B	181	VAL	3.7
1	E	233	THR	3.7
1	A	54	LYS	3.7
1	C	56	LEU	3.7
1	C	290	SER	3.7
1	C	237	SER	3.6
1	D	50	THR	3.6
1	B	237	SER	3.6
1	C	166	ALA	3.6
1	D	96	PRO	3.6
1	E	154	ASN	3.6
1	D	322	PHE	3.6
1	C	291	GLN	3.5
1	C	161	TRP	3.5
1	C	184	ASN	3.5
1	C	200	ASN	3.5
1	A	233	THR	3.5
1	B	165	LYS	3.5
1	D	162	ILE	3.5
1	D	93	MET	3.5
1	A	257	PRO	3.5
1	D	176	ASP	3.5
1	B	233	THR	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	179	SER	3.5
1	A	254	PRO	3.4
1	A	237	SER	3.4
1	A	49	LYS	3.4
1	B	258	TYR	3.4
1	D	99	ARG	3.4
1	B	182	GLN	3.4
1	D	95	PHE	3.4
1	D	70	GLY	3.4
1	D	44	THR	3.4
1	D	319	PHE	3.4
1	B	43	TRP	3.4
1	B	116	PHE	3.4
1	B	256	THR	3.4
1	D	291	GLN	3.4
1	D	256	THR	3.3
1	E	147	VAL	3.3
1	A	42	GLN	3.3
1	D	177	HIS	3.3
1	D	116	PHE	3.3
1	E	49	LYS	3.3
1	A	256	THR	3.2
1	A	117	ARG	3.2
1	D	101	ILE	3.2
1	E	161	TRP	3.2
1	A	258	TYR	3.2
1	C	122	ASP	3.2
1	C	47	PRO	3.1
1	E	153	ASP	3.1
1	B	151	ASN	3.1
1	B	166	ALA	3.1
1	B	87	THR	3.1
1	D	233	THR	3.1
1	C	145	ILE	3.1
1	B	47	PRO	3.1
1	D	92	LEU	3.1
1	B	121	PHE	3.1
1	C	55	PRO	3.0
1	D	180	SER	3.0
1	C	240	ILE	3.0
1	C	157	ILE	3.0
1	C	45	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	295	ALA	3.0
1	C	49	LYS	3.0
1	D	259	MET	2.9
1	A	50	THR	2.9
1	A	138	GLN	2.9
1	B	48	ARG	2.9
1	B	259	MET	2.9
1	D	164	GLY	2.9
1	B	254	PRO	2.9
1	B	202	PHE	2.9
1	E	155	GLU	2.9
1	D	43	TRP	2.9
1	D	51	PRO	2.9
1	C	233	THR	2.8
1	A	240	ILE	2.8
1	B	240	ILE	2.8
1	D	160	TRP	2.8
1	B	54	LYS	2.8
1	B	322	PHE	2.8
1	B	52	GLY	2.8
1	D	237	SER	2.8
1	D	65	ARG	2.8
1	D	157	ILE	2.8
1	A	95	PHE	2.8
1	D	296	ALA	2.8
1	C	158	ASP	2.8
1	C	286	LEU	2.7
1	C	126	PHE	2.7
1	B	290	SER	2.7
1	C	23	ILE	2.7
1	D	54	LYS	2.7
1	B	135	TYR	2.7
1	C	117	ARG	2.7
1	E	134	SER	2.7
1	A	51	PRO	2.7
1	B	45	GLY	2.7
1	A	321	PHE	2.7
1	D	100	VAL	2.7
1	C	116	PHE	2.7
1	A	289	GLU	2.7
1	B	161	TRP	2.6
1	B	150	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	163	ARG	2.6
1	E	159	GLU	2.6
1	C	202	PHE	2.6
1	B	241	ALA	2.6
1	B	154	ASN	2.6
1	D	197	ALA	2.6
1	C	177	HIS	2.6
1	E	58	VAL	2.6
1	A	116	PHE	2.6
1	A	99	ARG	2.6
1	B	155	GLU	2.6
1	B	185	GLN	2.6
1	A	285	TYR	2.6
1	A	161	TRP	2.6
1	A	93	MET	2.5
1	C	196	ASP	2.5
1	C	238	THR	2.5
1	C	31	GLN	2.5
1	A	303	ARG	2.5
1	C	241	ALA	2.5
1	C	87	THR	2.5
1	B	13	ASP	2.5
1	D	321	PHE	2.5
1	C	285	TYR	2.5
1	A	13	ASP	2.5
1	E	319	PHE	2.5
1	E	45	GLY	2.5
1	D	292	PRO	2.5
1	B	120	PRO	2.5
1	C	119	PHE	2.5
1	D	152	ILE	2.5
1	D	295	ALA	2.4
1	C	52	GLY	2.4
1	B	42	GLN	2.4
1	B	84	SER	2.4
1	E	294	ARG	2.4
1	D	240	ILE	2.4
1	A	101	ILE	2.4
1	A	56	LEU	2.4
1	E	148	TYR	2.4
1	B	321	PHE	2.4
1	D	135	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	114	MET	2.4
1	E	52	GLY	2.4
1	D	42	GLN	2.4
1	B	99	ARG	2.4
1	D	97	ASP	2.4
1	C	134	SER	2.3
1	D	288	VAL	2.3
1	D	293	ALA	2.3
1	C	57	ILE	2.3
1	E	135	TYR	2.3
1	D	68	ASN	2.3
1	E	292	PRO	2.3
1	E	113	ASP	2.3
1	E	241	ALA	2.3
1	C	256	THR	2.3
1	C	135	TYR	2.3
1	C	151	ASN	2.3
1	D	183	PRO	2.2
1	D	238	THR	2.2
1	B	291	GLN	2.2
1	B	257	PRO	2.2
1	E	181	VAL	2.2
1	D	74	PRO	2.2
1	E	285	TYR	2.2
1	C	118	LEU	2.2
1	D	84	SER	2.2
1	A	188	PHE	2.2
1	C	46	LYS	2.2
1	A	96	PRO	2.2
1	C	181	VAL	2.1
1	B	285	TYR	2.1
1	C	123	ARG	2.1
1	E	242	HIS	2.1
1	E	291	GLN	2.1
1	E	238	THR	2.1
1	B	94	LEU	2.1
1	C	84	SER	2.1
1	D	290	SER	2.1
1	A	236	VAL	2.1
1	C	88	GLY	2.1
1	D	236	VAL	2.1
1	D	242	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	198	VAL	2.1
1	C	288	VAL	2.1
1	B	162	ILE	2.1
1	A	139	GLN	2.1
1	A	224	TRP	2.1
1	E	254	PRO	2.1
1	D	257	PRO	2.1
1	A	88	GLY	2.0
1	B	118	LEU	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 carbohydrates 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.