



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 26, 2023 – 06:08 PM EDT

PDB ID : 8T7Z
Title : Crystal structure of alpha-glucosidase (yicI) from *Klebsiella aerogenes*
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2023-06-21
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.33
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) : 2.33

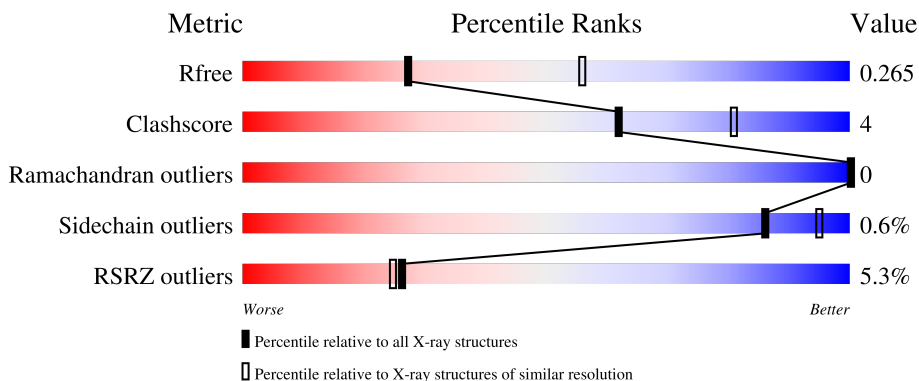
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	780	
1	B	780	
1	C	780	
1	D	780	
1	E	780	

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Mol	Chain	Length	Quality of chain
1	F	780	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a small red segment on the left labeled '4%', a large green segment in the middle labeled '86%', and a small yellow segment on the right labeled '10%'. A small grey dot is visible at the far right end of the bar.</p>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 36248 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 Alpha-glucosidase yicI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	754	6033	3865	1031	1109	28	0	0	0
1	B	758	6064	3881	1033	1122	28	0	0	0
1	C	758	6065	3881	1037	1119	28	0	0	0
1	D	751	6024	3859	1030	1107	28	0	0	0
1	E	753	6018	3857	1026	1107	28	0	0	0
1	F	752	6014	3853	1027	1106	28	0	0	0

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP A0A0H3FLJ3
A	2	ALA	-	expression tag	UNP A0A0H3FLJ3
A	3	HIS	-	expression tag	UNP A0A0H3FLJ3
A	4	HIS	-	expression tag	UNP A0A0H3FLJ3
A	5	HIS	-	expression tag	UNP A0A0H3FLJ3
A	6	HIS	-	expression tag	UNP A0A0H3FLJ3
A	7	HIS	-	expression tag	UNP A0A0H3FLJ3
A	8	HIS	-	expression tag	UNP A0A0H3FLJ3
A	81	GLN	HIS	engineered mutation	UNP A0A0H3FLJ3
A	87	SER	GLY	engineered mutation	UNP A0A0H3FLJ3
A	702	GLY	ASP	engineered mutation	UNP A0A0H3FLJ3
A	705	ASN	ARG	engineered mutation	UNP A0A0H3FLJ3
A	707	THR	ALA	engineered mutation	UNP A0A0H3FLJ3
A	708	VAL	LEU	engineered mutation	UNP A0A0H3FLJ3
A	718	ALA	VAL	engineered mutation	UNP A0A0H3FLJ3
B	1	MET	-	initiating methionine	UNP A0A0H3FLJ3
B	2	ALA	-	expression tag	UNP A0A0H3FLJ3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	3	HIS	-	expression tag	UNP A0A0H3FLJ3
B	4	HIS	-	expression tag	UNP A0A0H3FLJ3
B	5	HIS	-	expression tag	UNP A0A0H3FLJ3
B	6	HIS	-	expression tag	UNP A0A0H3FLJ3
B	7	HIS	-	expression tag	UNP A0A0H3FLJ3
B	8	HIS	-	expression tag	UNP A0A0H3FLJ3
B	81	GLN	HIS	engineered mutation	UNP A0A0H3FLJ3
B	87	SER	GLY	engineered mutation	UNP A0A0H3FLJ3
B	702	GLY	ASP	engineered mutation	UNP A0A0H3FLJ3
B	705	ASN	ARG	engineered mutation	UNP A0A0H3FLJ3
B	707	THR	ALA	engineered mutation	UNP A0A0H3FLJ3
B	708	VAL	LEU	engineered mutation	UNP A0A0H3FLJ3
B	718	ALA	VAL	engineered mutation	UNP A0A0H3FLJ3
C	1	MET	-	initiating methionine	UNP A0A0H3FLJ3
C	2	ALA	-	expression tag	UNP A0A0H3FLJ3
C	3	HIS	-	expression tag	UNP A0A0H3FLJ3
C	4	HIS	-	expression tag	UNP A0A0H3FLJ3
C	5	HIS	-	expression tag	UNP A0A0H3FLJ3
C	6	HIS	-	expression tag	UNP A0A0H3FLJ3
C	7	HIS	-	expression tag	UNP A0A0H3FLJ3
C	8	HIS	-	expression tag	UNP A0A0H3FLJ3
C	81	GLN	HIS	engineered mutation	UNP A0A0H3FLJ3
C	87	SER	GLY	engineered mutation	UNP A0A0H3FLJ3
C	702	GLY	ASP	engineered mutation	UNP A0A0H3FLJ3
C	705	ASN	ARG	engineered mutation	UNP A0A0H3FLJ3
C	707	THR	ALA	engineered mutation	UNP A0A0H3FLJ3
C	708	VAL	LEU	engineered mutation	UNP A0A0H3FLJ3
C	718	ALA	VAL	engineered mutation	UNP A0A0H3FLJ3
D	1	MET	-	initiating methionine	UNP A0A0H3FLJ3
D	2	ALA	-	expression tag	UNP A0A0H3FLJ3
D	3	HIS	-	expression tag	UNP A0A0H3FLJ3
D	4	HIS	-	expression tag	UNP A0A0H3FLJ3
D	5	HIS	-	expression tag	UNP A0A0H3FLJ3
D	6	HIS	-	expression tag	UNP A0A0H3FLJ3
D	7	HIS	-	expression tag	UNP A0A0H3FLJ3
D	8	HIS	-	expression tag	UNP A0A0H3FLJ3
D	81	GLN	HIS	engineered mutation	UNP A0A0H3FLJ3
D	87	SER	GLY	engineered mutation	UNP A0A0H3FLJ3
D	702	GLY	ASP	engineered mutation	UNP A0A0H3FLJ3
D	705	ASN	ARG	engineered mutation	UNP A0A0H3FLJ3
D	707	THR	ALA	engineered mutation	UNP A0A0H3FLJ3
D	708	VAL	LEU	engineered mutation	UNP A0A0H3FLJ3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	718	ALA	VAL	engineered mutation	UNP A0A0H3FLJ3
E	1	MET	-	initiating methionine	UNP A0A0H3FLJ3
E	2	ALA	-	expression tag	UNP A0A0H3FLJ3
E	3	HIS	-	expression tag	UNP A0A0H3FLJ3
E	4	HIS	-	expression tag	UNP A0A0H3FLJ3
E	5	HIS	-	expression tag	UNP A0A0H3FLJ3
E	6	HIS	-	expression tag	UNP A0A0H3FLJ3
E	7	HIS	-	expression tag	UNP A0A0H3FLJ3
E	8	HIS	-	expression tag	UNP A0A0H3FLJ3
E	81	GLN	HIS	engineered mutation	UNP A0A0H3FLJ3
E	87	SER	GLY	engineered mutation	UNP A0A0H3FLJ3
E	702	GLY	ASP	engineered mutation	UNP A0A0H3FLJ3
E	705	ASN	ARG	engineered mutation	UNP A0A0H3FLJ3
E	707	THR	ALA	engineered mutation	UNP A0A0H3FLJ3
E	708	VAL	LEU	engineered mutation	UNP A0A0H3FLJ3
E	718	ALA	VAL	engineered mutation	UNP A0A0H3FLJ3
F	1	MET	-	initiating methionine	UNP A0A0H3FLJ3
F	2	ALA	-	expression tag	UNP A0A0H3FLJ3
F	3	HIS	-	expression tag	UNP A0A0H3FLJ3
F	4	HIS	-	expression tag	UNP A0A0H3FLJ3
F	5	HIS	-	expression tag	UNP A0A0H3FLJ3
F	6	HIS	-	expression tag	UNP A0A0H3FLJ3
F	7	HIS	-	expression tag	UNP A0A0H3FLJ3
F	8	HIS	-	expression tag	UNP A0A0H3FLJ3
F	81	GLN	HIS	engineered mutation	UNP A0A0H3FLJ3
F	87	SER	GLY	engineered mutation	UNP A0A0H3FLJ3
F	702	GLY	ASP	engineered mutation	UNP A0A0H3FLJ3
F	705	ASN	ARG	engineered mutation	UNP A0A0H3FLJ3
F	707	THR	ALA	engineered mutation	UNP A0A0H3FLJ3
F	708	VAL	LEU	engineered mutation	UNP A0A0H3FLJ3
F	718	ALA	VAL	engineered mutation	UNP A0A0H3FLJ3

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total O 3 3	0	0
2	B	6	Total O 6 6	0	0
2	C	8	Total O 8 8	0	0
2	D	5	Total O 5 5	0	0

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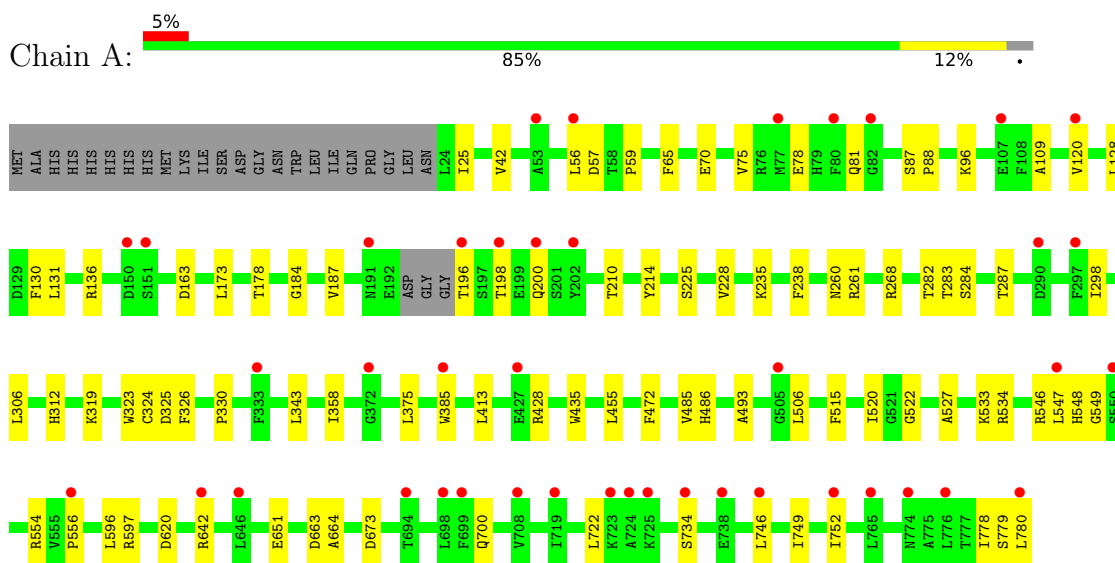
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	E	4	Total O 4 4	0	0
2	F	4	Total O 4 4	0	0

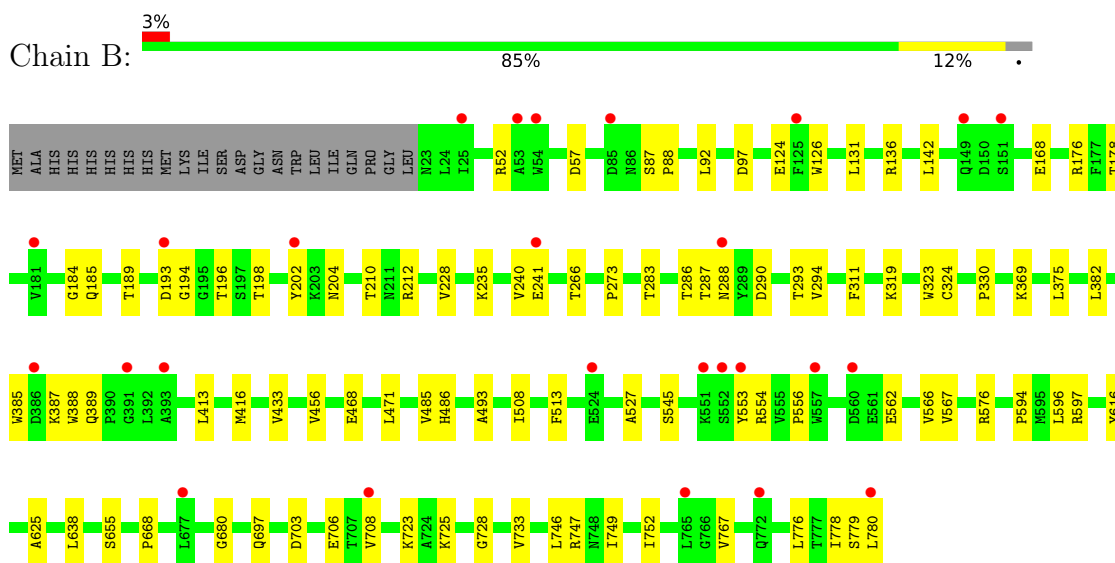
3 Residue-property plots [i](#)

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.

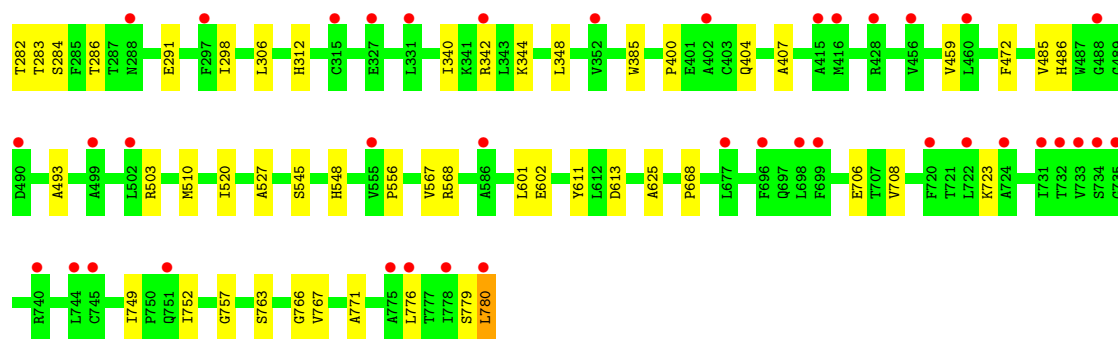
- Molecule 1: Alpha-glucosidase yicI



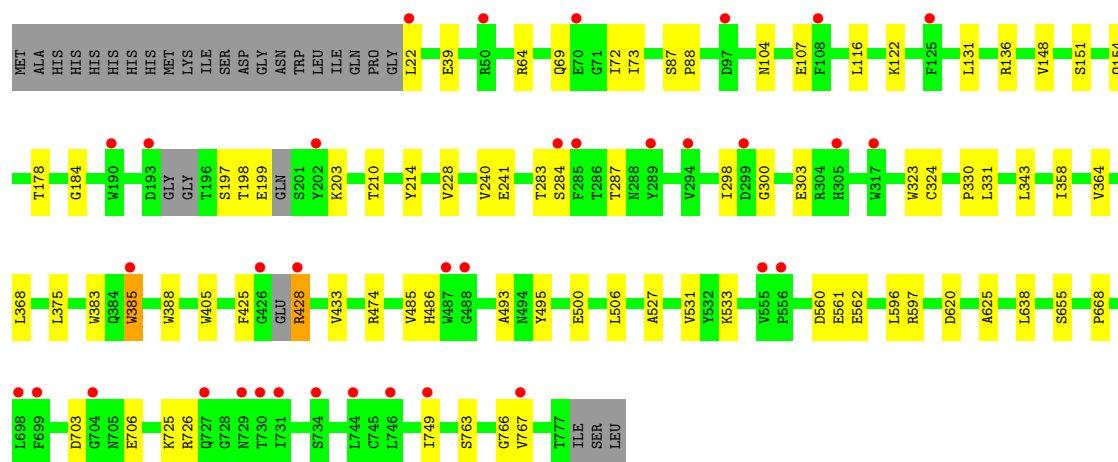
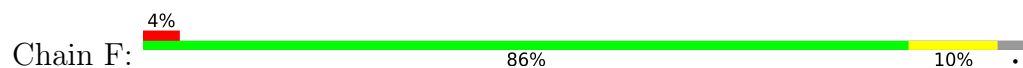
- Molecule 1: Alpha-glucosidase yicI



- Molecule 1: Alpha-glucosidase yicI



• Molecule 1: Alpha-glucosidase yicI



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.79Å 208.51Å 280.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.25 – 2.70 49.55 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.25-2.70) 100.0 (49.55-2.70)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 2.69Å)	Xtrriage
Refinement program	PHENIX (1.21rc1_4933: ???)	Depositor
R, R_{free}	0.229 , 0.263 0.233 , 0.265	Depositor DCC
R_{free} test set	6572 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	47.2	Xtrriage
Anisotropy	0.627	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	36248	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.28	0/6211	0.52	0/8441
1	B	0.28	0/6243	0.52	0/8485
1	C	0.28	0/6244	0.53	0/8486
1	D	0.27	0/6202	0.52	0/8427
1	E	0.29	0/6195	0.52	0/8419
1	F	0.27	0/6190	0.51	0/8412
All	All	0.28	0/37285	0.52	0/50670

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	6033	0	5756	70	0
1	B	6064	0	5776	56	0
1	C	6065	0	5783	61	0
1	D	6024	0	5751	37	0
1	E	6018	0	5730	51	0
1	F	6014	0	5729	47	0
2	A	3	0	0	0	0
2	B	6	0	0	1	0
2	C	8	0	0	0	0
2	D	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	4	0	0	0	0
2	F	4	0	0	0	0
All	All	36248	0	34525	309	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.

The worst 5 of 309 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:F:210:THR:HG22	1:F:214:TYR:O	1.86	0.76
1:D:168:GLU:OE2	1:D:210:THR:HG21	1.87	0.74
1:B:168:GLU:OE2	1:B:210:THR:HG21	1.88	0.73
1:A:210:THR:HG22	1:A:214:TYR:O	1.89	0.73
1:A:642:ARG:CZ	1:A:651:GLU:HB3	2.19	0.72

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	750/780 (96%)	727 (97%)	23 (3%)	0	100	100
1	B	756/780 (97%)	728 (96%)	28 (4%)	0	100	100
1	C	756/780 (97%)	731 (97%)	25 (3%)	0	100	100
1	D	747/780 (96%)	718 (96%)	29 (4%)	0	100	100
1	E	747/780 (96%)	722 (97%)	25 (3%)	0	100	100
1	F	744/780 (95%)	718 (96%)	26 (4%)	0	100	100
All	All	4500/4680 (96%)	4344 (96%)	156 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	634/660 (96%)	630 (99%)	4 (1%)	86	95
1	B	638/660 (97%)	632 (99%)	6 (1%)	78	92
1	C	638/660 (97%)	635 (100%)	3 (0%)	88	96
1	D	634/660 (96%)	630 (99%)	4 (1%)	86	95
1	E	631/660 (96%)	627 (99%)	4 (1%)	86	95
1	F	632/660 (96%)	630 (100%)	2 (0%)	92	98
All	All	3807/3960 (96%)	3784 (99%)	23 (1%)	86	95

5 of 23 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	176	ARG
1	E	203	LYS
1	D	780	LEU
1	E	385	TRP
1	B	176	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	29	GLN

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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	754/780 (96%)	0.47	42 (5%) 24 23	35, 51, 76, 107	0
1	B	758/780 (97%)	0.33	26 (3%) 45 45	35, 49, 82, 126	0
1	C	758/780 (97%)	0.47	55 (7%) 15 13	38, 51, 78, 129	0
1	D	751/780 (96%)	0.30	24 (3%) 47 48	35, 48, 68, 102	0
1	E	753/780 (96%)	0.64	60 (7%) 12 10	38, 56, 80, 123	0
1	F	752/780 (96%)	0.37	35 (4%) 31 30	36, 50, 76, 106	0
All	All	4526/4680 (96%)	0.43	242 (5%) 26 25	35, 51, 77, 129	0

The worst 5 of 242 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	753	ALA	6.5
1	C	776	LEU	6.5
1	E	731	ILE	6.4
1	E	733	VAL	5.9
1	E	402	ALA	5.5

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.