Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 09:25 AM EDT

PDB ID	:	5E9A
Title	:	Crystal structure analysis of the cold-adamped beta-galactosidase from Rahnella sp. R3
Authors	:	Zhang, Y.Z.; Fan, Y.T.
Deposited on	:	2015-10-14
Resolution	:	2.56 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

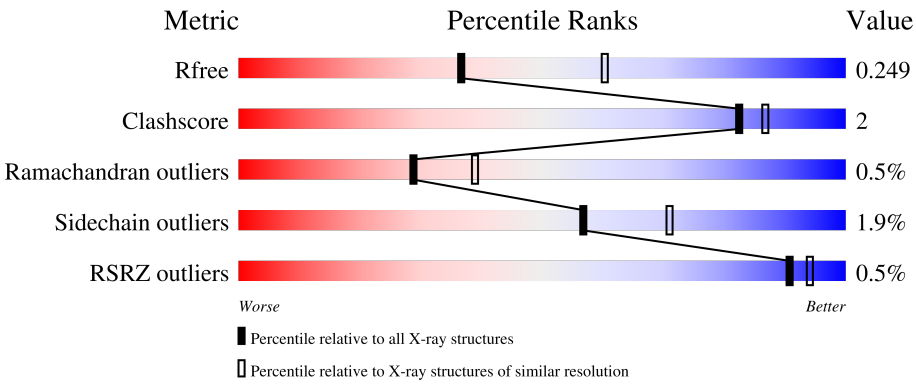
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.56 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	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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	712	86% 9% • 5%
1	B	712	% 88% 8% • •
1	C	712	% 90% 6% • •
1	D	712	88% 8% •
1	E	712	% 86% 9% • •

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Mol	Chain	Length	Quality of chain
1	F	712	88% 7% • •

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
3	ACT	F	702	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 32981 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 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	679	5387	3420	945	993	29	0	0	0
1	B	684	5424	3443	952	1000	29	0	0	0
1	C	684	5424	3443	952	1000	29	0	0	0
1	D	684	5424	3443	952	1000	29	0	0	0
1	E	684	5432	3448	955	1000	29	0	1	0
1	F	684	5432	3448	955	1000	29	0	1	0

There are 150 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	initiating methionine	UNP A0A0B4U8I5
A	-23	ASN	-	expression tag	UNP A0A0B4U8I5
A	-22	HIS	-	expression tag	UNP A0A0B4U8I5
A	-21	LYS	-	expression tag	UNP A0A0B4U8I5
A	-20	VAL	-	expression tag	UNP A0A0B4U8I5
A	-19	HIS	-	expression tag	UNP A0A0B4U8I5
A	-18	HIS	-	expression tag	UNP A0A0B4U8I5
A	-17	HIS	-	expression tag	UNP A0A0B4U8I5
A	-16	HIS	-	expression tag	UNP A0A0B4U8I5
A	-15	HIS	-	expression tag	UNP A0A0B4U8I5
A	-14	HIS	-	expression tag	UNP A0A0B4U8I5
A	-13	ILE	-	expression tag	UNP A0A0B4U8I5
A	-12	GLU	-	expression tag	UNP A0A0B4U8I5
A	-11	GLY	-	expression tag	UNP A0A0B4U8I5
A	-10	ARG	-	expression tag	UNP A0A0B4U8I5
A	-9	HIS	-	expression tag	UNP A0A0B4U8I5
A	-8	MET	-	expression tag	UNP A0A0B4U8I5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLU	-	expression tag	UNP A0A0B4U8I5
A	-6	LEU	-	expression tag	UNP A0A0B4U8I5
A	-5	GLY	-	expression tag	UNP A0A0B4U8I5
A	-4	THR	-	expression tag	UNP A0A0B4U8I5
A	-3	LEU	-	expression tag	UNP A0A0B4U8I5
A	-2	GLU	-	expression tag	UNP A0A0B4U8I5
A	-1	GLY	-	expression tag	UNP A0A0B4U8I5
A	0	SER	-	expression tag	UNP A0A0B4U8I5
B	-24	MET	-	initiating methionine	UNP A0A0B4U8I5
B	-23	ASN	-	expression tag	UNP A0A0B4U8I5
B	-22	HIS	-	expression tag	UNP A0A0B4U8I5
B	-21	LYS	-	expression tag	UNP A0A0B4U8I5
B	-20	VAL	-	expression tag	UNP A0A0B4U8I5
B	-19	HIS	-	expression tag	UNP A0A0B4U8I5
B	-18	HIS	-	expression tag	UNP A0A0B4U8I5
B	-17	HIS	-	expression tag	UNP A0A0B4U8I5
B	-16	HIS	-	expression tag	UNP A0A0B4U8I5
B	-15	HIS	-	expression tag	UNP A0A0B4U8I5
B	-14	HIS	-	expression tag	UNP A0A0B4U8I5
B	-13	ILE	-	expression tag	UNP A0A0B4U8I5
B	-12	GLU	-	expression tag	UNP A0A0B4U8I5
B	-11	GLY	-	expression tag	UNP A0A0B4U8I5
B	-10	ARG	-	expression tag	UNP A0A0B4U8I5
B	-9	HIS	-	expression tag	UNP A0A0B4U8I5
B	-8	MET	-	expression tag	UNP A0A0B4U8I5
B	-7	GLU	-	expression tag	UNP A0A0B4U8I5
B	-6	LEU	-	expression tag	UNP A0A0B4U8I5
B	-5	GLY	-	expression tag	UNP A0A0B4U8I5
B	-4	THR	-	expression tag	UNP A0A0B4U8I5
B	-3	LEU	-	expression tag	UNP A0A0B4U8I5
B	-2	GLU	-	expression tag	UNP A0A0B4U8I5
B	-1	GLY	-	expression tag	UNP A0A0B4U8I5
B	0	SER	-	expression tag	UNP A0A0B4U8I5
C	-24	MET	-	initiating methionine	UNP A0A0B4U8I5
C	-23	ASN	-	expression tag	UNP A0A0B4U8I5
C	-22	HIS	-	expression tag	UNP A0A0B4U8I5
C	-21	LYS	-	expression tag	UNP A0A0B4U8I5
C	-20	VAL	-	expression tag	UNP A0A0B4U8I5
C	-19	HIS	-	expression tag	UNP A0A0B4U8I5
C	-18	HIS	-	expression tag	UNP A0A0B4U8I5
C	-17	HIS	-	expression tag	UNP A0A0B4U8I5
C	-16	HIS	-	expression tag	UNP A0A0B4U8I5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-15	HIS	-	expression tag	UNP A0A0B4U8I5
C	-14	HIS	-	expression tag	UNP A0A0B4U8I5
C	-13	ILE	-	expression tag	UNP A0A0B4U8I5
C	-12	GLU	-	expression tag	UNP A0A0B4U8I5
C	-11	GLY	-	expression tag	UNP A0A0B4U8I5
C	-10	ARG	-	expression tag	UNP A0A0B4U8I5
C	-9	HIS	-	expression tag	UNP A0A0B4U8I5
C	-8	MET	-	expression tag	UNP A0A0B4U8I5
C	-7	GLU	-	expression tag	UNP A0A0B4U8I5
C	-6	LEU	-	expression tag	UNP A0A0B4U8I5
C	-5	GLY	-	expression tag	UNP A0A0B4U8I5
C	-4	THR	-	expression tag	UNP A0A0B4U8I5
C	-3	LEU	-	expression tag	UNP A0A0B4U8I5
C	-2	GLU	-	expression tag	UNP A0A0B4U8I5
C	-1	GLY	-	expression tag	UNP A0A0B4U8I5
C	0	SER	-	expression tag	UNP A0A0B4U8I5
D	-24	MET	-	initiating methionine	UNP A0A0B4U8I5
D	-23	ASN	-	expression tag	UNP A0A0B4U8I5
D	-22	HIS	-	expression tag	UNP A0A0B4U8I5
D	-21	LYS	-	expression tag	UNP A0A0B4U8I5
D	-20	VAL	-	expression tag	UNP A0A0B4U8I5
D	-19	HIS	-	expression tag	UNP A0A0B4U8I5
D	-18	HIS	-	expression tag	UNP A0A0B4U8I5
D	-17	HIS	-	expression tag	UNP A0A0B4U8I5
D	-16	HIS	-	expression tag	UNP A0A0B4U8I5
D	-15	HIS	-	expression tag	UNP A0A0B4U8I5
D	-14	HIS	-	expression tag	UNP A0A0B4U8I5
D	-13	ILE	-	expression tag	UNP A0A0B4U8I5
D	-12	GLU	-	expression tag	UNP A0A0B4U8I5
D	-11	GLY	-	expression tag	UNP A0A0B4U8I5
D	-10	ARG	-	expression tag	UNP A0A0B4U8I5
D	-9	HIS	-	expression tag	UNP A0A0B4U8I5
D	-8	MET	-	expression tag	UNP A0A0B4U8I5
D	-7	GLU	-	expression tag	UNP A0A0B4U8I5
D	-6	LEU	-	expression tag	UNP A0A0B4U8I5
D	-5	GLY	-	expression tag	UNP A0A0B4U8I5
D	-4	THR	-	expression tag	UNP A0A0B4U8I5
D	-3	LEU	-	expression tag	UNP A0A0B4U8I5
D	-2	GLU	-	expression tag	UNP A0A0B4U8I5
D	-1	GLY	-	expression tag	UNP A0A0B4U8I5
D	0	SER	-	expression tag	UNP A0A0B4U8I5
E	-24	MET	-	initiating methionine	UNP A0A0B4U8I5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-23	ASN	-	expression tag	UNP A0A0B4U8I5
E	-22	HIS	-	expression tag	UNP A0A0B4U8I5
E	-21	LYS	-	expression tag	UNP A0A0B4U8I5
E	-20	VAL	-	expression tag	UNP A0A0B4U8I5
E	-19	HIS	-	expression tag	UNP A0A0B4U8I5
E	-18	HIS	-	expression tag	UNP A0A0B4U8I5
E	-17	HIS	-	expression tag	UNP A0A0B4U8I5
E	-16	HIS	-	expression tag	UNP A0A0B4U8I5
E	-15	HIS	-	expression tag	UNP A0A0B4U8I5
E	-14	HIS	-	expression tag	UNP A0A0B4U8I5
E	-13	ILE	-	expression tag	UNP A0A0B4U8I5
E	-12	GLU	-	expression tag	UNP A0A0B4U8I5
E	-11	GLY	-	expression tag	UNP A0A0B4U8I5
E	-10	ARG	-	expression tag	UNP A0A0B4U8I5
E	-9	HIS	-	expression tag	UNP A0A0B4U8I5
E	-8	MET	-	expression tag	UNP A0A0B4U8I5
E	-7	GLU	-	expression tag	UNP A0A0B4U8I5
E	-6	LEU	-	expression tag	UNP A0A0B4U8I5
E	-5	GLY	-	expression tag	UNP A0A0B4U8I5
E	-4	THR	-	expression tag	UNP A0A0B4U8I5
E	-3	LEU	-	expression tag	UNP A0A0B4U8I5
E	-2	GLU	-	expression tag	UNP A0A0B4U8I5
E	-1	GLY	-	expression tag	UNP A0A0B4U8I5
E	0	SER	-	expression tag	UNP A0A0B4U8I5
F	-24	MET	-	initiating methionine	UNP A0A0B4U8I5
F	-23	ASN	-	expression tag	UNP A0A0B4U8I5
F	-22	HIS	-	expression tag	UNP A0A0B4U8I5
F	-21	LYS	-	expression tag	UNP A0A0B4U8I5
F	-20	VAL	-	expression tag	UNP A0A0B4U8I5
F	-19	HIS	-	expression tag	UNP A0A0B4U8I5
F	-18	HIS	-	expression tag	UNP A0A0B4U8I5
F	-17	HIS	-	expression tag	UNP A0A0B4U8I5
F	-16	HIS	-	expression tag	UNP A0A0B4U8I5
F	-15	HIS	-	expression tag	UNP A0A0B4U8I5
F	-14	HIS	-	expression tag	UNP A0A0B4U8I5
F	-13	ILE	-	expression tag	UNP A0A0B4U8I5
F	-12	GLU	-	expression tag	UNP A0A0B4U8I5
F	-11	GLY	-	expression tag	UNP A0A0B4U8I5
F	-10	ARG	-	expression tag	UNP A0A0B4U8I5
F	-9	HIS	-	expression tag	UNP A0A0B4U8I5
F	-8	MET	-	expression tag	UNP A0A0B4U8I5
F	-7	GLU	-	expression tag	UNP A0A0B4U8I5

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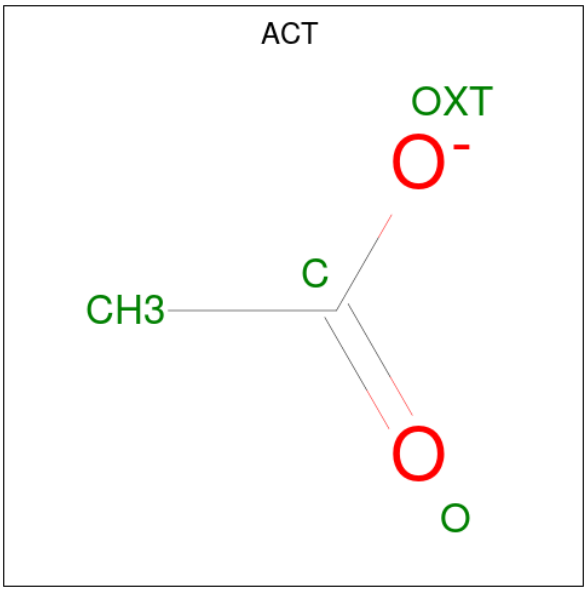
Chain	Residue	Modelled	Actual	Comment	Reference
F	-6	LEU	-	expression tag	UNP A0A0B4U8I5
F	-5	GLY	-	expression tag	UNP A0A0B4U8I5
F	-4	THR	-	expression tag	UNP A0A0B4U8I5
F	-3	LEU	-	expression tag	UNP A0A0B4U8I5
F	-2	GLU	-	expression tag	UNP A0A0B4U8I5
F	-1	GLY	-	expression tag	UNP A0A0B4U8I5
F	0	SER	-	expression tag	UNP A0A0B4U8I5

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 4 is water.

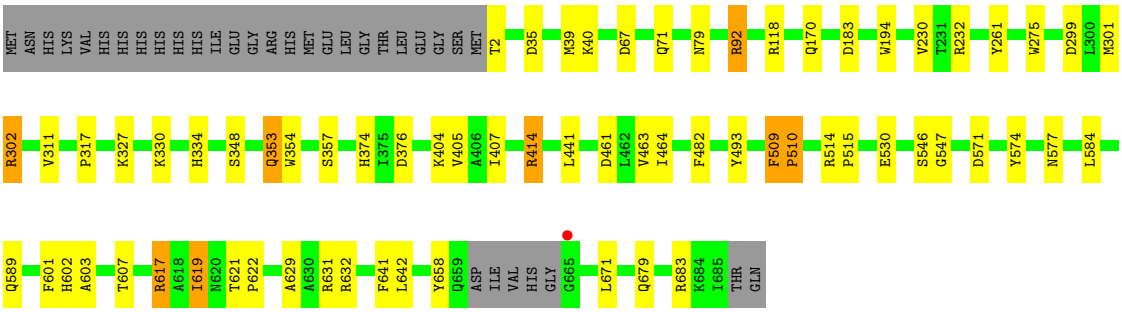
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	105	105	105	0	0
4	B	56	56	56	0	0
4	C	66	66	66	0	0
4	D	99	99	99	0	0
4	E	48	48	48	0	0
4	F	54	54	54	0	0

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.

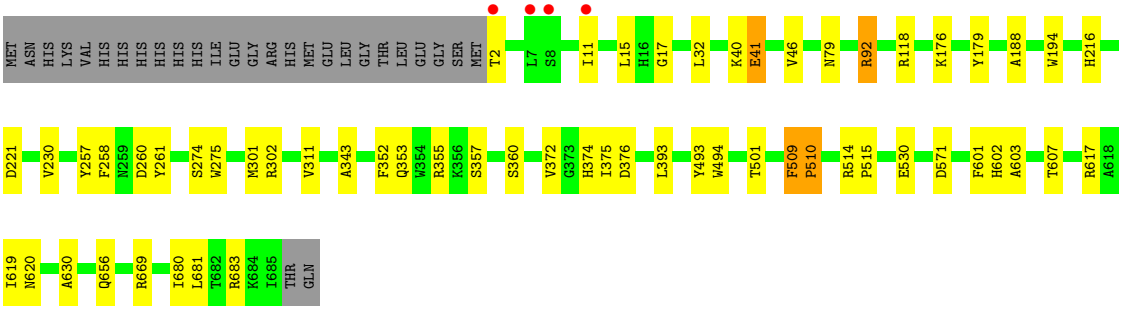
• Molecule 1: Beta-galactosidase

Chain A:



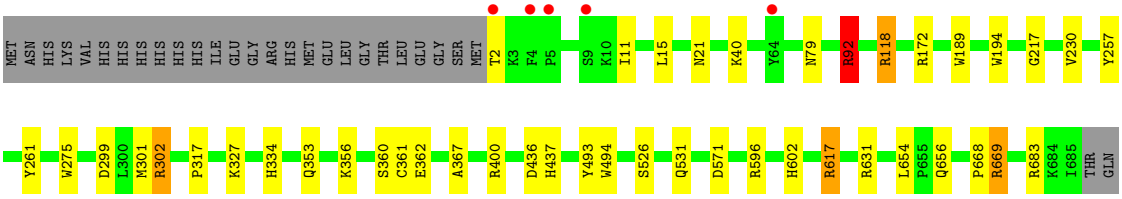
• Molecule 1: Beta-galactosidase

Chain B:



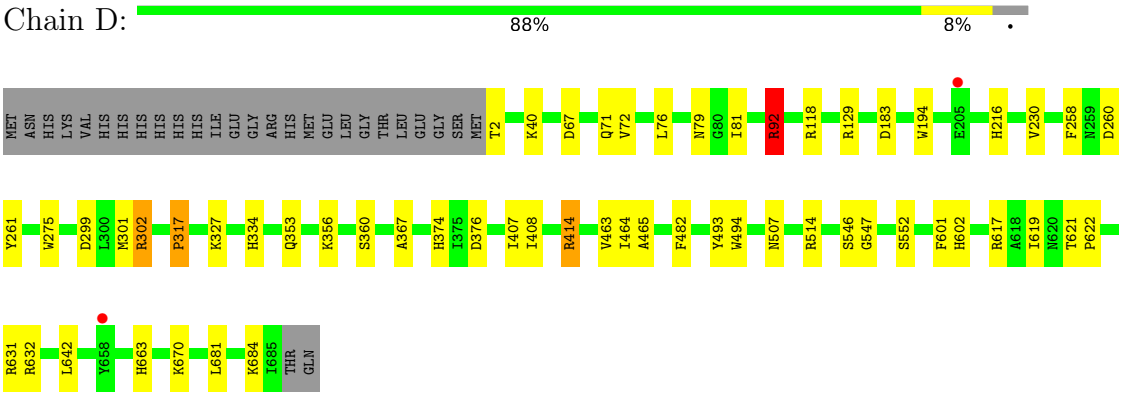
• Molecule 1: Beta-galactosidase

Chain C:



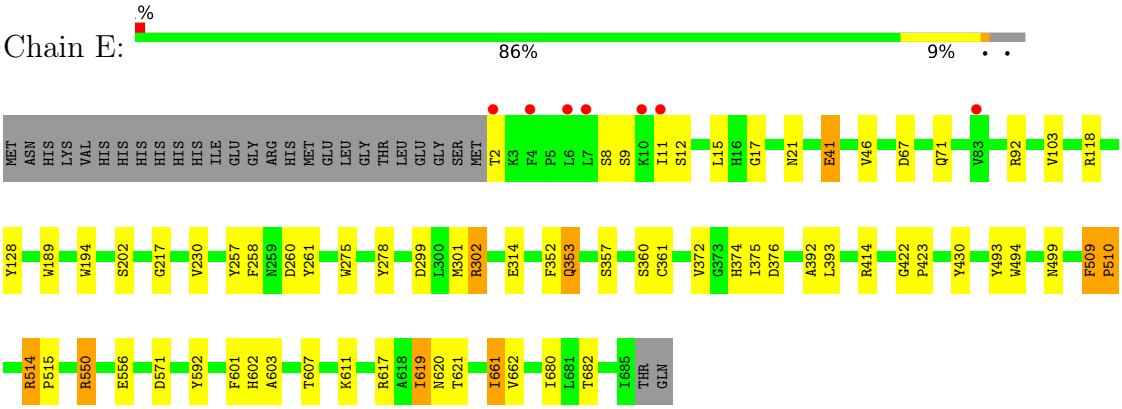
• Molecule 1: Beta-galactosidase

Chain D:



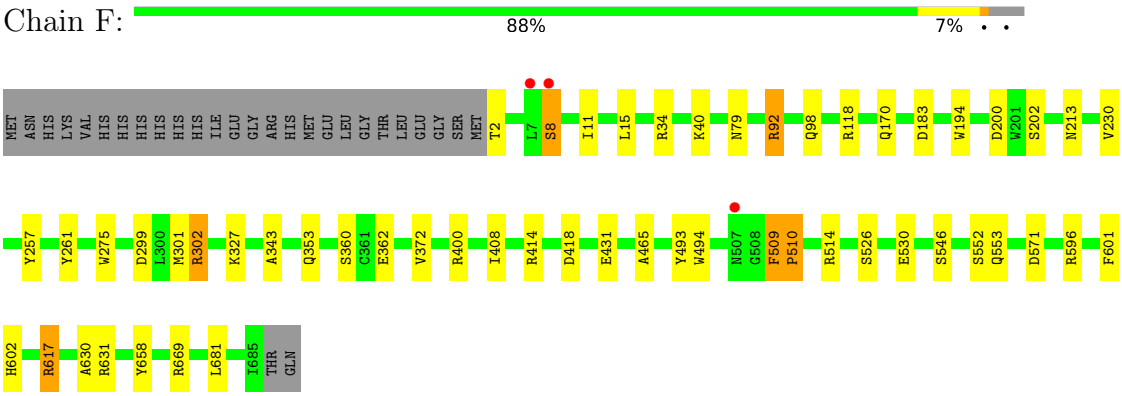
• Molecule 1: Beta-galactosidase

Chain E:



• Molecule 1: Beta-galactosidase

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	146.43Å 106.83Å 164.24Å 90.00° 109.00° 90.00°	Depositor
Resolution (Å)	138.45 – 2.56 138.45 – 2.56	Depositor EDS
% Data completeness (in resolution range)	98.5 (138.45-2.56) 98.5 (138.45-2.56)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$< I/\sigma(I) >$ [1]	2.08 (at 2.55Å)	Xtriage
Refinement program	REFMAC 5.8.0131	Depositor
R, R_{free}	0.199 , 0.247 0.204 , 0.249	Depositor DCC
R_{free} test set	2020 reflections (1.33%)	wwPDB-VP
Wilson B-factor (Å^2)	29.6	Xtriage
Anisotropy	0.099	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 39.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	32981	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.29% 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: ACT, 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	0.69	0/5540	0.85	11/7537 (0.1%)
1	B	0.65	0/5579	0.81	9/7592 (0.1%)
1	C	0.65	0/5579	0.82	13/7592 (0.2%)
1	D	0.69	0/5579	0.83	8/7592 (0.1%)
1	E	0.63	0/5590	0.81	9/7607 (0.1%)
1	F	0.63	0/5590	0.81	10/7606 (0.1%)
All	All	0.66	0/33457	0.82	60/45526 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	D	0	1
1	E	0	2
1	F	0	1
All	All	0	6

There are no bond length outliers.

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	92	ARG	NE-CZ-NH1	13.27	126.94	120.30
1	D	92	ARG	NE-CZ-NH1	12.79	126.69	120.30
1	A	509	PHE	C-N-CD	-10.72	97.03	120.60
1	C	92	ARG	NE-CZ-NH2	-9.66	115.47	120.30
1	F	509	PHE	C-N-CD	-9.58	99.52	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	302	ARG	NE-CZ-NH1	9.40	125.00	120.30
1	D	92	ARG	NE-CZ-NH2	-8.99	115.80	120.30
1	F	92	ARG	NE-CZ-NH1	8.68	124.64	120.30
1	E	550	ARG	NE-CZ-NH2	8.63	124.61	120.30
1	E	509	PHE	C-N-CD	-8.51	101.88	120.60
1	A	302	ARG	NE-CZ-NH2	-8.15	116.22	120.30
1	D	302	ARG	NE-CZ-NH1	7.92	124.26	120.30
1	B	509	PHE	C-N-CD	-7.92	103.18	120.60
1	C	617	ARG	NE-CZ-NH1	7.54	124.07	120.30
1	C	302	ARG	NE-CZ-NH1	7.43	124.02	120.30
1	C	302	ARG	NE-CZ-NH2	-7.11	116.75	120.30
1	F	302	ARG	NE-CZ-NH1	6.97	123.78	120.30
1	F	92	ARG	NE-CZ-NH2	-6.91	116.84	120.30
1	F	400	ARG	NE-CZ-NH1	6.70	123.65	120.30
1	E	302	ARG	NE-CZ-NH1	6.61	123.61	120.30
1	D	302	ARG	NE-CZ-NH2	-6.48	117.06	120.30
1	E	611	LYS	CD-CE-NZ	6.34	126.28	111.70
1	F	669	ARG	CG-CD-NE	-6.08	99.03	111.80
1	E	661	ILE	CB-CA-C	-6.02	99.56	111.60
1	B	302	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	C	400	ARG	NE-CZ-NH1	5.85	123.22	120.30
1	B	352	PHE	N-CA-CB	5.84	121.11	110.60
1	B	302	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	C	617	ARG	NE-CZ-NH2	-5.79	117.40	120.30
1	D	414	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	C	596	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	B	221	ASP	CB-CG-OD2	-5.77	113.11	118.30
1	A	414	ARG	NE-CZ-NH1	5.69	123.14	120.30
1	A	617	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	A	631	ARG	NE-CZ-NH1	5.51	123.05	120.30
1	B	355	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	E	352	PHE	N-CA-CB	5.49	120.48	110.60
1	F	631	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	E	619	ILE	CB-CA-C	-5.48	100.64	111.60
1	A	617	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	C	118	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	C	631	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	B	352	PHE	O-C-N	5.31	131.20	122.70
1	B	619	ILE	CB-CA-C	-5.31	100.99	111.60
1	B	352	PHE	C-N-CA	5.29	134.93	121.70
1	C	436	ASP	CB-CG-OD2	-5.29	113.54	118.30
1	D	92	ARG	CD-NE-CZ	5.29	131.01	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	632	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	E	302	ARG	NE-CZ-NH2	-5.27	117.67	120.30
1	F	418	ASP	CB-CG-OD1	5.26	123.03	118.30
1	A	232	ARG	NE-CZ-NH1	5.25	122.92	120.30
1	D	631	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	A	619	ILE	CB-CA-C	-5.18	101.24	111.60
1	F	617	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	A	632	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	E	514	ARG	NE-CZ-NH1	5.11	122.85	120.30
1	C	92	ARG	CD-NE-CZ	5.06	130.68	123.60
1	F	302	ARG	NE-CZ-NH2	-5.05	117.78	120.30
1	C	172	ARG	NE-CZ-NH1	5.03	122.82	120.30
1	A	509	PHE	C-N-CA	5.01	143.03	122.00

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	509	PHE	Peptide
1	B	509	PHE	Peptide
1	D	670	LYS	Peptide
1	E	422	GLY	Peptide
1	E	509	PHE	Peptide
1	F	509	PHE	Peptide

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	5387	0	5151	38	0
1	B	5424	0	5186	26	0
1	C	5424	0	5186	22	0
1	D	5424	0	5186	28	0
1	E	5432	0	5199	33	0
1	F	5432	0	5199	22	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
3	C	4	0	3	1	0
3	D	4	0	3	0	0
3	E	4	0	3	1	0
3	F	4	0	3	2	0
4	A	105	0	0	5	0
4	B	56	0	0	1	0
4	C	66	0	0	2	0
4	D	99	0	0	2	0
4	E	48	0	0	2	0
4	F	54	0	0	3	0
All	All	32981	0	31125	159	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 2.

All (159) 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)
3:E:702:ACT:H2	4:E:825:HOH:O	1.81	0.80
1:E:661:ILE:HD11	1:E:682:THR:HB	1.66	0.78
1:F:414:ARG:HD2	4:F:808:HOH:O	1.90	0.71
1:D:92:ARG:HG2	1:D:92:ARG:HH11	1.57	0.69
1:E:230:VAL:HG21	1:E:261:TYR:CZ	2.28	0.69
1:C:362:GLU:OE2	3:C:702:ACT:H2	1.93	0.68
1:F:230:VAL:HG21	1:F:261:TYR:CZ	2.28	0.67
1:B:656:GLN:OE1	1:B:683:ARG:NH1	2.28	0.66
1:C:230:VAL:HG21	1:C:261:TYR:CZ	2.31	0.65
1:E:661:ILE:HD11	1:E:682:THR:CB	2.26	0.65
1:B:230:VAL:HG21	1:B:261:TYR:CZ	2.34	0.63
1:D:299:ASP:OD1	1:D:302:ARG:NH2	2.31	0.63
1:A:230:VAL:HG21	1:A:261:TYR:CZ	2.35	0.61
1:F:275:TRP:CD1	1:F:301:MET:HG3	2.36	0.61
1:F:327:LYS:NZ	4:F:801:HOH:O	2.33	0.60
1:D:275:TRP:CD1	1:D:301:MET:HG3	2.36	0.60
1:A:619:ILE:HD11	1:A:642:LEU:HD21	1.81	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:275:TRP:CD1	1:C:301:MET:HG3	2.38	0.59
1:D:230:VAL:HG21	1:D:261:TYR:CZ	2.38	0.59
1:E:353:GLN:NE2	4:E:801:HOH:O	2.36	0.59
1:A:353:GLN:NE2	4:A:801:HOH:O	2.35	0.59
1:C:92:ARG:HG2	1:C:92:ARG:HH11	1.68	0.58
3:F:702:ACT:H1	4:F:842:HOH:O	2.02	0.58
1:F:362:GLU:OE2	3:F:702:ACT:H2	2.03	0.57
1:D:507:ASN:O	1:D:514:ARG:NH1	2.37	0.57
1:E:11:ILE:HG21	1:E:15:LEU:HD21	1.86	0.56
1:A:92:ARG:HH11	1:A:92:ARG:N	2.05	0.55
1:C:437:HIS:HA	1:C:602:HIS:HD2	1.72	0.55
1:D:414:ARG:HD2	4:D:818:HOH:O	2.06	0.55
1:A:601:PHE:HD2	1:A:602:HIS:HD2	1.55	0.54
1:C:40:LYS:HD2	1:C:79:ASN:HB3	1.90	0.54
1:A:275:TRP:CD1	1:A:301:MET:HG3	2.43	0.53
1:A:299:ASP:OD1	1:A:302:ARG:NH2	2.42	0.53
1:E:275:TRP:CD1	1:E:301:MET:HG3	2.44	0.52
1:F:510:PRO:HD2	1:F:514:ARG:HB2	1.90	0.52
1:A:619:ILE:CD1	1:A:642:LEU:HD21	2.40	0.52
1:A:374:HIS:HD2	1:A:376:ASP:H	1.58	0.52
1:D:601:PHE:HD2	1:D:602:HIS:HD2	1.58	0.52
1:E:194:TRP:CE2	1:F:118:ARG:HD3	2.44	0.51
1:E:202:SER:OG	1:F:98:GLN:NE2	2.43	0.51
1:C:654:LEU:O	1:C:669:ARG:HG3	2.09	0.51
1:B:11:ILE:HG21	1:B:15:LEU:HD21	1.92	0.51
1:B:601:PHE:HD2	1:B:602:HIS:HD2	1.59	0.51
1:B:603:ALA:O	1:B:607:THR:HG23	2.11	0.51
1:E:601:PHE:HD2	1:E:602:HIS:HD2	1.58	0.51
1:A:67:ASP:O	1:A:71:GLN:HG2	2.11	0.50
1:B:374:HIS:HD2	1:B:376:ASP:H	1.58	0.50
1:F:658:TYR:CD2	1:F:681:LEU:HD23	2.47	0.49
1:A:546:SER:OG	1:A:547:GLY:N	2.45	0.49
1:E:499:ASN:OD1	1:E:499:ASN:C	2.51	0.49
1:A:330:LYS:NZ	4:A:802:HOH:O	2.43	0.49
1:C:654:LEU:HD12	1:C:668:PRO:O	2.12	0.49
1:B:92:ARG:O	1:B:92:ARG:HD2	2.12	0.49
1:A:510:PRO:HD2	1:A:514:ARG:HB2	1.94	0.49
1:E:17:GLY:HA3	1:E:46:VAL:O	2.12	0.49
1:B:681:LEU:HD12	1:B:681:LEU:N	2.28	0.48
1:D:619:ILE:HD11	1:D:642:LEU:HD21	1.95	0.48
1:C:656:GLN:OE1	1:C:683:ARG:CZ	2.62	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:327:LYS:HG2	1:F:526:SER:O	2.14	0.48
1:E:392:ALA:HB3	1:E:662:VAL:HG12	1.95	0.47
1:B:393:LEU:HD11	1:B:680:ILE:HD13	1.95	0.47
1:B:194:TRP:CE2	1:C:118:ARG:HD3	2.50	0.47
1:F:40:LYS:HD2	1:F:79:ASN:HB3	1.96	0.47
1:A:603:ALA:O	1:A:607:THR:HG23	2.14	0.47
1:D:546:SER:OG	1:D:547:GLY:N	2.48	0.47
1:A:414:ARG:HD2	4:A:820:HOH:O	2.14	0.47
1:D:356:LYS:HG3	1:D:367:ALA:HB3	1.97	0.47
1:A:35:ASP:O	1:A:39:MET:HG3	2.15	0.47
1:B:374:HIS:HD2	1:B:376:ASP:N	2.13	0.47
1:B:17:GLY:HA3	1:B:46:VAL:O	2.14	0.46
1:B:275:TRP:CD1	1:B:301:MET:HG3	2.49	0.46
1:B:514:ARG:HB3	1:B:515:PRO:HD3	1.96	0.46
1:D:408:ILE:HD12	1:D:465:ALA:HB2	1.96	0.46
1:A:327:LYS:HG2	1:C:526:SER:O	2.16	0.46
1:A:629:ALA:HA	1:A:641:PHE:O	2.15	0.46
1:C:327:LYS:NZ	4:C:804:HOH:O	2.47	0.46
1:D:194:TRP:CE2	1:E:118:ARG:HD3	2.51	0.46
1:A:40:LYS:HD2	1:A:79:ASN:HB3	1.98	0.46
1:D:40:LYS:HD3	1:D:79:ASN:HB3	1.97	0.46
1:E:510:PRO:HD2	1:E:514:ARG:HB2	1.98	0.46
1:D:129:ARG:HG3	4:D:832:HOH:O	2.15	0.46
1:F:200:ASP:OD2	1:F:202:SER:OG	2.21	0.46
1:A:407:ILE:HG12	1:A:464:ILE:HB	1.98	0.45
1:D:407:ILE:HG12	1:D:464:ILE:HB	1.98	0.45
1:F:601:PHE:HD2	1:F:602:HIS:HD2	1.64	0.45
1:B:393:LEU:HD11	1:B:680:ILE:CD1	2.47	0.45
1:A:317:PRO:HA	1:A:334:HIS:CE1	2.51	0.45
1:D:317:PRO:HA	1:D:334:HIS:CE1	2.51	0.45
1:E:603:ALA:O	1:E:607:THR:HG23	2.17	0.45
1:A:194:TRP:CE2	1:B:118:ARG:HD3	2.51	0.45
1:E:67:ASP:O	1:E:71:GLN:HG2	2.17	0.45
1:D:118:ARG:HD3	1:F:194:TRP:CE2	2.51	0.45
1:F:408:ILE:HD12	1:F:465:ALA:HB2	1.99	0.45
1:E:41:GLU:HG2	1:E:375:ILE:HD11	1.99	0.45
1:B:32:LEU:HB3	4:B:850:HOH:O	2.17	0.44
1:B:176:LYS:O	1:B:179:TYR:O	2.35	0.44
1:D:463:VAL:HG23	1:D:482:PHE:CZ	2.52	0.44
1:B:258:PHE:CE2	1:B:260:ASP:HB2	2.52	0.44
1:B:274:SER:HA	1:B:311:VAL:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:577:ASN:HB3	4:A:862:HOH:O	2.16	0.44
1:E:661:ILE:CD1	1:E:682:THR:HG21	2.47	0.44
1:C:92:ARG:HH11	1:C:92:ARG:CG	2.30	0.44
1:D:67:ASP:O	1:D:71:GLN:HG2	2.17	0.44
1:A:39:MET:HA	1:A:354:TRP:CH2	2.53	0.43
1:D:374:HIS:HD2	1:D:376:ASP:H	1.66	0.43
1:D:258:PHE:CE2	1:D:260:ASP:HB2	2.53	0.43
1:A:353:GLN:HG2	4:A:892:HOH:O	2.18	0.43
1:F:213:ASN:OD1	1:F:213:ASN:N	2.51	0.43
1:A:514:ARG:HB3	1:A:515:PRO:HD3	2.00	0.43
1:A:584:LEU:HD12	1:A:589:GLN:HG2	2.00	0.43
1:A:311:VAL:HG13	1:A:348:SER:HB3	2.00	0.43
1:C:356:LYS:HG3	1:C:367:ALA:HB3	2.01	0.43
1:B:216:HIS:NE2	1:C:361:CYS:HA	2.34	0.43
1:A:658:TYR:HD1	1:A:683:ARG:HB2	1.83	0.43
1:E:278:TYR:OH	1:E:314:GLU:HB3	2.19	0.43
1:E:103:VAL:HG12	1:E:128:TYR:HB2	2.01	0.43
1:E:550:ARG:HH21	1:E:550:ARG:HG2	1.84	0.43
1:A:327:LYS:HE2	1:C:531:GLN:OE1	2.19	0.42
1:E:556:GLU:OE1	1:E:592:TYR:OH	2.22	0.42
1:F:11:ILE:HG21	1:F:15:LEU:HD21	2.00	0.42
1:E:299:ASP:OD1	1:E:302:ARG:NH2	2.52	0.42
1:F:414:ARG:NH1	1:F:431:GLU:OE2	2.53	0.42
1:A:327:LYS:NZ	4:C:805:HOH:O	2.51	0.42
1:B:510:PRO:HD2	1:B:514:ARG:HB2	2.01	0.42
1:F:553:GLN:OE1	1:F:596:ARG:NH1	2.44	0.42
1:E:374:HIS:HD2	1:E:376:ASP:N	2.17	0.42
1:A:118:ARG:HD3	1:C:194:TRP:CE2	2.55	0.42
1:B:188:ALA:HB1	1:B:501:THR:HG21	2.02	0.42
1:E:189:TRP:CD1	1:E:217:GLY:HA3	2.55	0.42
1:A:404:LYS:HD3	1:A:461:ASP:OD2	2.19	0.41
1:A:621:THR:OG1	1:A:622:PRO:HD2	2.20	0.41
1:A:671:LEU:HD11	1:A:679:GLN:OE1	2.20	0.41
1:E:374:HIS:HD2	1:E:376:ASP:H	1.68	0.41
1:E:510:PRO:HD2	1:E:514:ARG:CB	2.50	0.41
1:C:189:TRP:CD1	1:C:217:GLY:HA3	2.55	0.41
1:A:405:VAL:HG11	1:A:441:LEU:HD13	2.01	0.41
1:B:41:GLU:HG2	1:B:375:ILE:HD11	2.02	0.41
1:E:514:ARG:HB3	1:E:515:PRO:HD3	2.02	0.41
1:D:621:THR:OG1	1:D:622:PRO:HD2	2.20	0.41
1:E:258:PHE:CE2	1:E:260:ASP:HB2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:619:ILE:CG2	1:E:621:THR:HG22	2.51	0.41
1:A:463:VAL:HG23	1:A:482:PHE:CZ	2.56	0.41
1:B:40:LYS:HD2	1:B:79:ASN:HB3	2.03	0.41
1:C:11:ILE:HG21	1:C:15:LEU:HD21	2.02	0.41
1:C:317:PRO:HA	1:C:334:HIS:CE1	2.56	0.41
1:D:76:LEU:HD22	1:D:81:ILE:HG21	2.02	0.41
1:D:216:HIS:NE2	1:E:361:CYS:HA	2.36	0.41
1:D:619:ILE:CD1	1:D:642:LEU:HD21	2.50	0.41
1:D:681:LEU:N	1:D:681:LEU:HD12	2.36	0.41
1:F:299:ASP:OD1	1:F:302:ARG:NH2	2.54	0.41
1:C:299:ASP:OD1	1:C:302:ARG:NH2	2.54	0.40
1:A:574:TYR:O	1:A:577:ASN:HB2	2.21	0.40
1:B:343:ALA:HB1	1:B:630:ALA:HB3	2.02	0.40
1:D:642:LEU:HD23	1:D:642:LEU:HA	1.92	0.40
1:F:343:ALA:HB1	1:F:630:ALA:HB3	2.02	0.40
1:F:8:SER:HB3	1:F:11:ILE:HG12	2.03	0.40
1:E:393:LEU:HD11	1:E:680:ILE:HD13	2.03	0.40
1:E:414:ARG:HD3	1:E:430:TYR:HD2	1.85	0.40
1:C:92:ARG:HG2	1:C:92:ARG:NH1	2.34	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	675/712 (95%)	650 (96%)	22 (3%)	3 (0%)	34 45
1	B	682/712 (96%)	652 (96%)	26 (4%)	4 (1%)	25 33
1	C	682/712 (96%)	654 (96%)	25 (4%)	3 (0%)	34 45
1	D	682/712 (96%)	654 (96%)	25 (4%)	3 (0%)	34 45
1	E	683/712 (96%)	649 (95%)	29 (4%)	5 (1%)	22 29

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	683/712 (96%)	651 (95%)	28 (4%)	4 (1%)	25	33
All	All	4087/4272 (96%)	3910 (96%)	155 (4%)	22 (0%)	29	39

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	510	PRO
1	B	510	PRO
1	D	663	HIS
1	E	493	TYR
1	E	510	PRO
1	F	510	PRO
1	C	493	TYR
1	D	493	TYR
1	A	493	TYR
1	B	353	GLN
1	D	353	GLN
1	E	257	TYR
1	E	353	GLN
1	F	257	TYR
1	F	493	TYR
1	A	353	GLN
1	B	493	TYR
1	C	353	GLN
1	C	257	TYR
1	E	423	PRO
1	F	353	GLN
1	B	257	TYR

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	563/592 (95%)	555 (99%)	8 (1%)	67	78
1	B	567/592 (96%)	555 (98%)	12 (2%)	53	67

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	567/592 (96%)	559 (99%)	8 (1%)	67	78
1	D	567/592 (96%)	557 (98%)	10 (2%)	59	73
1	E	568/592 (96%)	554 (98%)	14 (2%)	47	61
1	F	568/592 (96%)	554 (98%)	14 (2%)	47	61
All	All	3400/3552 (96%)	3334 (98%)	66 (2%)	57	71

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	92	ARG
1	A	170	GLN
1	A	183	ASP
1	A	357	SER
1	A	530	GLU
1	A	571	ASP
1	A	617	ARG
1	B	2	THR
1	B	41	GLU
1	B	92	ARG
1	B	357	SER
1	B	360	SER
1	B	372	VAL
1	B	494	TRP
1	B	530	GLU
1	B	571	ASP
1	B	617	ARG
1	B	620	ASN
1	B	669	ARG
1	C	2	THR
1	C	21	ASN
1	C	92	ARG
1	C	360	SER
1	C	494	TRP
1	C	571	ASP
1	C	617	ARG
1	C	669	ARG
1	D	2	THR
1	D	72	VAL
1	D	92	ARG
1	D	183	ASP

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Mol	Chain	Res	Type
1	D	317	PRO
1	D	360	SER
1	D	494	TRP
1	D	552	SER
1	D	617	ARG
1	D	684	LYS
1	E	2	THR
1	E	8	SER
1	E	9	SER
1	E	12	SER
1	E	21	ASN
1	E	41	GLU
1	E	92	ARG
1	E	357	SER
1	E	360	SER
1	E	372	VAL
1	E	494	TRP
1	E	571	ASP
1	E	617	ARG
1	E	620	ASN
1	F	2	THR
1	F	8	SER
1	F	34	ARG
1	F	92	ARG
1	F	170	GLN
1	F	183	ASP
1	F	360	SER
1	F	372	VAL
1	F	494	TRP
1	F	530	GLU
1	F	546	SER
1	F	552	SER
1	F	571	ASP
1	F	617	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	374	HIS
1	B	374	HIS
1	C	374	HIS
1	C	602	HIS

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Mol	Chain	Res	Type
1	D	374	HIS
1	E	374	HIS
1	F	374	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 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	ACT	F	702	-	3,3,3	0.81	0	3,3,3	0.40	0
3	ACT	A	702	-	3,3,3	0.83	0	3,3,3	0.58	0
3	ACT	E	702	-	3,3,3	0.63	0	3,3,3	1.40	0
3	ACT	C	702	-	3,3,3	0.81	0	3,3,3	0.22	0
3	ACT	B	702	-	3,3,3	0.79	0	3,3,3	0.91	0
3	ACT	D	702	-	3,3,3	0.41	0	3,3,3	1.91	2 (66%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	702	ACT	OXT-C-CH3	2.43	125.24	115.18
3	D	702	ACT	O-C-CH3	-2.21	113.72	122.33

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	702	ACT	2	0
3	E	702	ACT	1	0
3	C	702	ACT	1	0

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	679/712 (95%)	-0.19	1 (0%)	95	98	15, 25, 44, 81	0
1	B	684/712 (96%)	-0.12	4 (0%)	89	93	19, 30, 49, 82	0
1	C	684/712 (96%)	-0.05	5 (0%)	87	91	19, 32, 52, 80	0
1	D	684/712 (96%)	-0.17	2 (0%)	94	97	16, 25, 45, 84	0
1	E	684/712 (96%)	0.02	7 (1%)	82	87	21, 34, 52, 87	0
1	F	684/712 (96%)	0.04	3 (0%)	92	96	21, 34, 56, 81	0
All	All	4099/4272 (95%)	-0.08	22 (0%)	91	94	15, 30, 51, 87	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	11	ILE	4.4
1	E	7	LEU	4.0
1	E	2	THR	3.8
1	B	7	LEU	3.7
1	E	6	LEU	3.2
1	B	2	THR	3.1
1	C	2	THR	3.0
1	E	4	PHE	2.9
1	B	11	ILE	2.7
1	F	507	ASN	2.6
1	E	10	LYS	2.6
1	A	665	GLY	2.5
1	C	9	SER	2.5
1	D	658	TYR	2.5
1	F	7	LEU	2.3
1	E	83	VAL	2.2
1	D	205	GLU	2.2
1	C	4	PHE	2.1
1	F	8	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	64	TYR	2.1
1	B	8	SER	2.1
1	C	5	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
3	ACT	D	702	4/4	0.84	0.18	34,37,40,45	0
3	ACT	B	702	4/4	0.85	0.16	35,38,40,43	0
3	ACT	C	702	4/4	0.88	0.17	41,42,44,49	0
3	ACT	A	702	4/4	0.88	0.20	35,37,39,43	0
3	ACT	F	702	4/4	0.90	0.13	41,43,44,45	0
3	ACT	E	702	4/4	0.96	0.13	30,32,32,34	0
2	ZN	A	701	1/1	0.99	0.10	24,24,24,24	0
2	ZN	B	701	1/1	0.99	0.08	32,32,32,32	0
2	ZN	C	701	1/1	0.99	0.09	34,34,34,34	0
2	ZN	D	701	1/1	0.99	0.10	27,27,27,27	0
2	ZN	E	701	1/1	0.99	0.09	37,37,37,37	0
2	ZN	F	701	1/1	0.99	0.09	36,36,36,36	0

6.5 Other polymers (i)

There are no such residues in this entry.