



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2024 – 04:27 AM EST

PDB ID : 3OB8  
Title : Structure of the beta-galactosidase from *Kluyveromyces lactis* in complex with galactose  
Authors : Fernandez-Leiro, R.; Pereira-Rodriguez, A.; Becerra, M.; Gonzalez-Siso, I.; Cerdan, M.E.; Sanz-Aparicio, J.  
Deposited on : 2010-08-06  
Resolution : 2.80 Å(reported)

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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>.

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The following versions of software and data (see [references ⓘ](#)) 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 : 2.36  
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) : 2.36

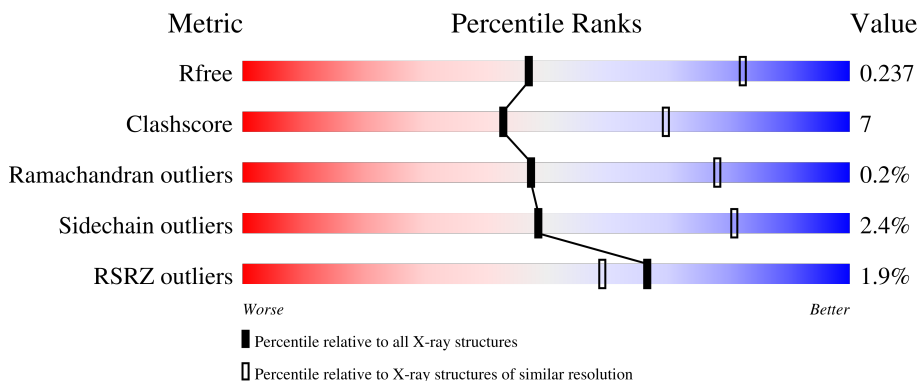
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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	1032	85% 13% ..
1	B	1032	85% 13% ..
1	C	1032	86% 12% ..
1	D	1032	85% 13% ..

## 2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 34419 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	1024	8325	5335	1388	1587	15	0	0	0
1	B	1024	8325	5335	1388	1587	15	0	0	0
1	D	1024	8325	5335	1388	1587	15	0	0	0
1	C	1024	8325	5335	1388	1587	15	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

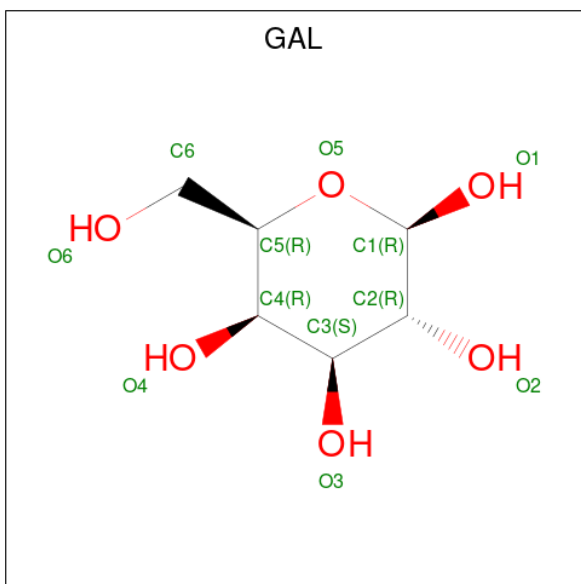
Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	ASP	-	expression tag	UNP P00723
A	-5	TYR	-	expression tag	UNP P00723
A	-4	LYS	-	expression tag	UNP P00723
A	-3	ASP	-	expression tag	UNP P00723
A	-2	ASP	-	expression tag	UNP P00723
A	-1	ASP	-	expression tag	UNP P00723
A	0	ASP	-	expression tag	UNP P00723
A	1	LYS	-	expression tag	UNP P00723
B	-6	ASP	-	expression tag	UNP P00723
B	-5	TYR	-	expression tag	UNP P00723
B	-4	LYS	-	expression tag	UNP P00723
B	-3	ASP	-	expression tag	UNP P00723
B	-2	ASP	-	expression tag	UNP P00723
B	-1	ASP	-	expression tag	UNP P00723
B	0	ASP	-	expression tag	UNP P00723
B	1	LYS	-	expression tag	UNP P00723
D	-6	ASP	-	expression tag	UNP P00723
D	-5	TYR	-	expression tag	UNP P00723
D	-4	LYS	-	expression tag	UNP P00723
D	-3	ASP	-	expression tag	UNP P00723
D	-2	ASP	-	expression tag	UNP P00723

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	ASP	-	expression tag	UNP P00723
D	0	ASP	-	expression tag	UNP P00723
D	1	LYS	-	expression tag	UNP P00723
C	-6	ASP	-	expression tag	UNP P00723
C	-5	TYR	-	expression tag	UNP P00723
C	-4	LYS	-	expression tag	UNP P00723
C	-3	ASP	-	expression tag	UNP P00723
C	-2	ASP	-	expression tag	UNP P00723
C	-1	ASP	-	expression tag	UNP P00723
C	0	ASP	-	expression tag	UNP P00723
C	1	LYS	-	expression tag	UNP P00723

- Molecule 2 is beta-D-galactopyranose (three-letter code: GAL) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf		
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		
2	D	1	Total	C	O	0	0
			12	6	6		
2	C	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0
3	D	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total Na 4 4	0	0
4	B	4	Total Na 4 4	0	0
4	D	4	Total Na 4 4	0	0
4	C	4	Total Na 4 4	0	0

- Molecule 5 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Mn 1 1	0	0
5	B	1	Total Mn 1 1	0	0
5	D	1	Total Mn 1 1	0	0
5	C	1	Total Mn 1 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	231	Total O 231 231	0	0
6	B	293	Total O 293 293	0	0
6	D	253	Total O 253 253	0	0

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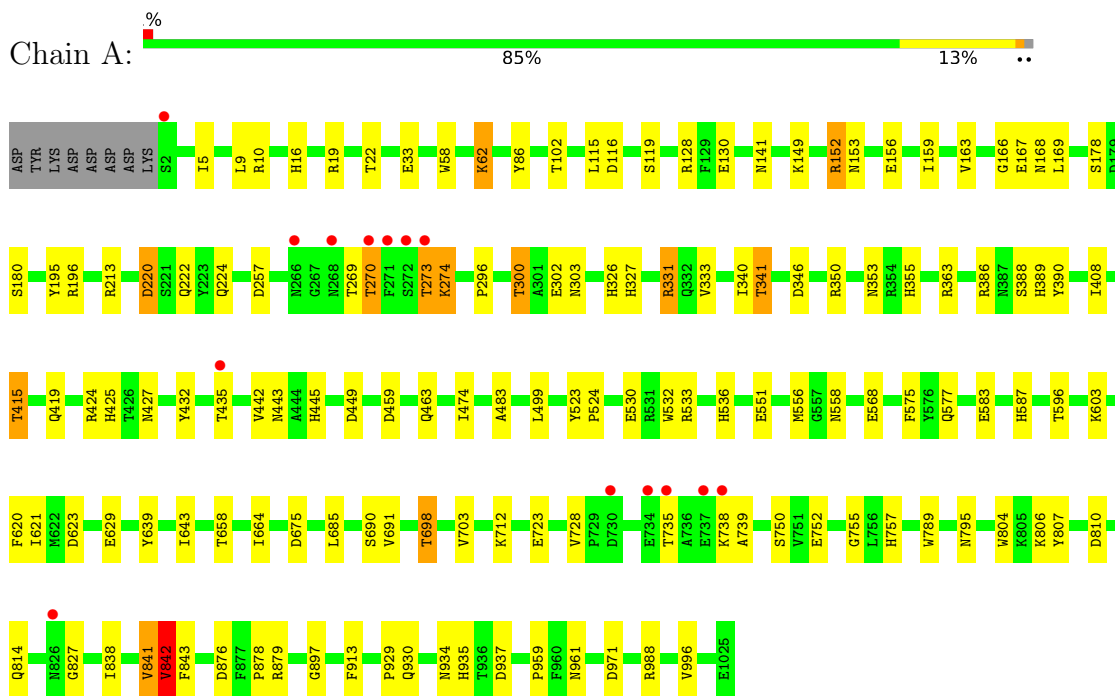
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
6	C	270	Total 270	O 270	0	0

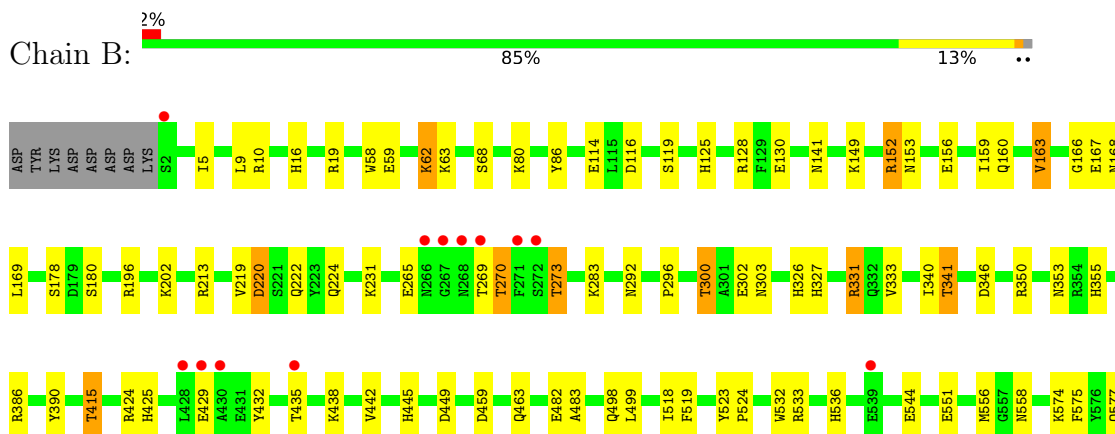
### 3 Residue-property plots [i](#)

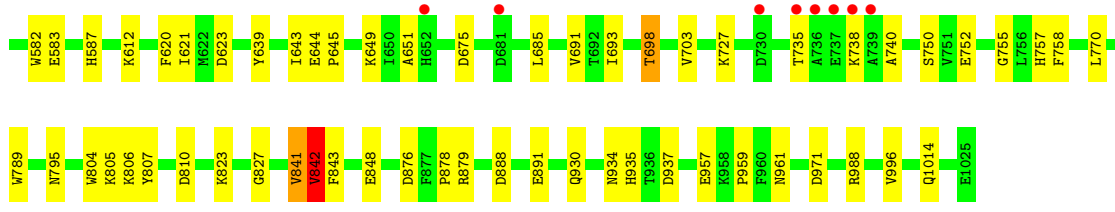
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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

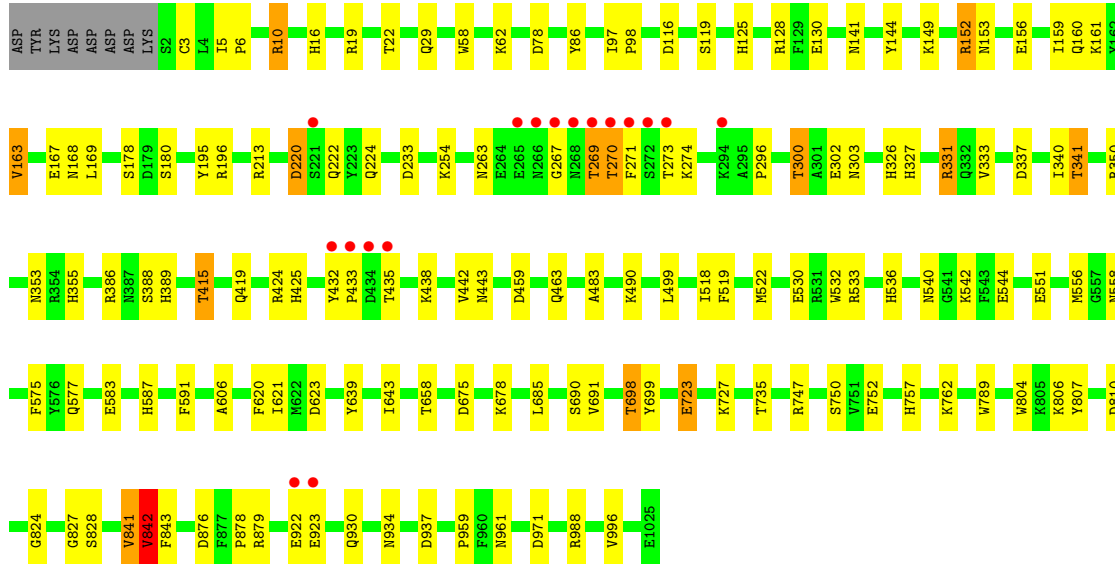
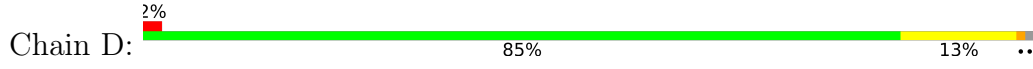


- Molecule 1: Beta-galactosidase

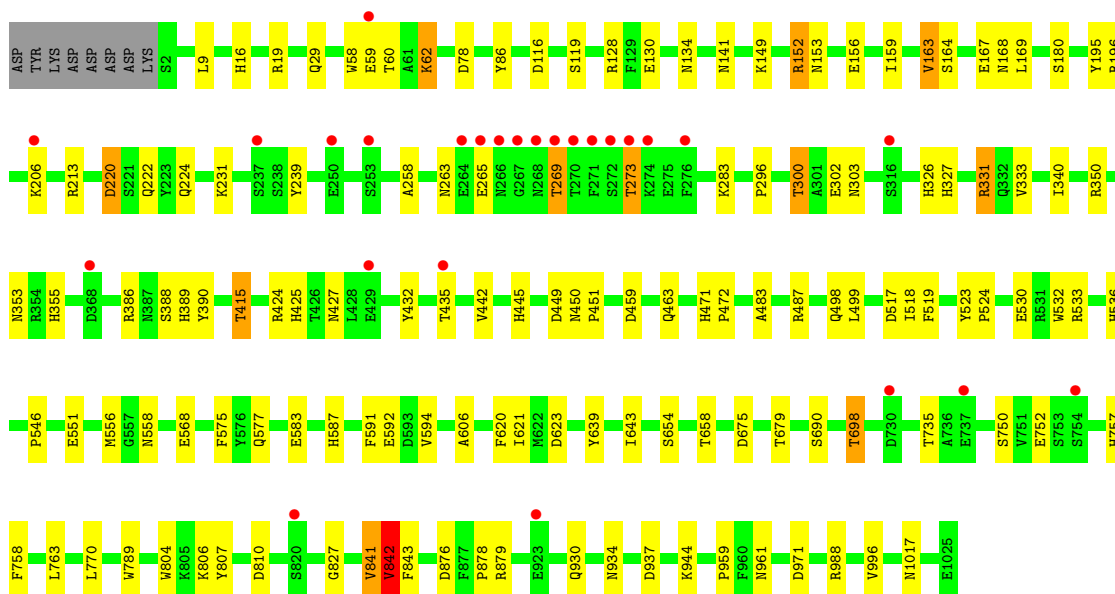
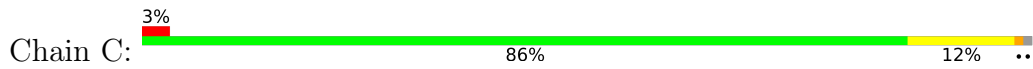




• Molecule 1: Beta-galactosidase



• Molecule 1: Beta-galactosidase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	140.38Å 153.45Å 217.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.30 – 2.80 49.30 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.30-2.80) 99.9 (49.30-2.80)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.35 (at 2.81Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.214 , 0.246 0.209 , 0.237	Depositor DCC
$R_{free}$ test set	5800 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.7	Xtrriage
Anisotropy	0.064	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 31.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	34419	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.14% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MN, GAL, NA, MG

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	5/8561 (0.1%)	0.56	1/11614 (0.0%)
1	B	0.50	5/8561 (0.1%)	0.56	0/11614
1	C	0.51	5/8561 (0.1%)	0.57	0/11614
1	D	0.51	5/8561 (0.1%)	0.56	0/11614
All	All	0.50	20/34244 (0.1%)	0.56	1/46456 (0.0%)

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	843	PHE	CD2-CE2	-5.80	1.27	1.39
1	B	843	PHE	CD2-CE2	-5.80	1.27	1.39
1	A	843	PHE	CD2-CE2	-5.79	1.27	1.39
1	D	843	PHE	CD2-CE2	-5.78	1.27	1.39
1	B	843	PHE	CD1-CE1	-5.48	1.28	1.39

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	346	ASP	CB-CG-OD2	5.07	122.87	118.30

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	8325	0	7972	119	0
1	B	8325	0	7972	126	0
1	C	8325	0	7972	112	1
1	D	8325	0	7972	120	1
2	A	12	0	12	1	0
2	B	12	0	12	1	0
2	C	12	0	11	1	0
2	D	12	0	11	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
4	C	4	0	0	0	0
4	D	4	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	A	231	0	0	19	0
6	B	293	0	0	34	0
6	C	270	0	0	28	0
6	D	253	0	0	32	0
All	All	34419	0	31934	446	1

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 7.

The worst 5 of 446 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:C:587:HIS:HB2	1:C:621:ILE:HG13	1.28	1.11
1:D:587:HIS:HB2	1:D:621:ILE:HG13	1.27	1.09
1:B:587:HIS:HB2	1:B:621:ILE:HG13	1.29	1.07
1:A:587:HIS:HB2	1:A:621:ILE:HG13	1.29	1.06
1:B:841:VAL:O	1:B:842:VAL:HG13	1.56	1.06

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:144:TYR:O	1:C:273:THR:OG1[3_545]	1.80	0.40

### 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	1022/1032 (99%)	971 (95%)	48 (5%)	3 (0%)	41 72
1	B	1022/1032 (99%)	971 (95%)	49 (5%)	2 (0%)	47 78
1	C	1022/1032 (99%)	973 (95%)	47 (5%)	2 (0%)	47 78
1	D	1022/1032 (99%)	971 (95%)	49 (5%)	2 (0%)	47 78
All	All	4088/4128 (99%)	3886 (95%)	193 (5%)	9 (0%)	47 78

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	842	VAL
1	B	842	VAL
1	D	842	VAL
1	C	842	VAL
1	A	62	LYS

#### 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	905/913 (99%)	883 (98%)	22 (2%)	49 81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	905/913 (99%)	883 (98%)	22 (2%)	49 81
1	C	905/913 (99%)	884 (98%)	21 (2%)	50 82
1	D	905/913 (99%)	882 (98%)	23 (2%)	47 80
All	All	3620/3652 (99%)	3532 (98%)	88 (2%)	49 81

5 of 88 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	442	VAL
1	C	167	GLU
1	D	620	PHE
1	D	971	ASP
1	C	300	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 98 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	387	ASN
1	D	934	ASN
1	D	419	GLN
1	D	536	HIS
1	C	125	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

Of 28 ligands modelled in this entry, 24 are monoatomic - leaving 4 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
2	GAL	B	2001	4	12,12,12	0.55	0	17,17,17	0.63	0
2	GAL	D	2001	4	12,12,12	0.47	0	17,17,17	1.05	1 (5%)
2	GAL	A	2001	4	12,12,12	0.55	0	17,17,17	0.98	1 (5%)
2	GAL	C	2001	4	12,12,12	0.53	0	17,17,17	1.25	1 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GAL	B	2001	4	-	2/2/22/22	0/1/1/1
2	GAL	D	2001	4	-	2/2/22/22	0/1/1/1
2	GAL	A	2001	4	-	2/2/22/22	0/1/1/1
2	GAL	C	2001	4	-	1/2/22/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2001	GAL	C1-O5-C5	3.18	119.67	113.66
2	D	2001	GAL	C1-O5-C5	2.51	118.40	113.66
2	A	2001	GAL	C1-O5-C5	2.45	118.30	113.66

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	2001	GAL	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	A	2001	GAL	O5-C5-C6-O6
2	D	2001	GAL	O5-C5-C6-O6
2	B	2001	GAL	C4-C5-C6-O6
2	A	2001	GAL	C4-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2001	GAL	1	0
2	D	2001	GAL	1	0
2	A	2001	GAL	1	0
2	C	2001	GAL	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	1024/1032 (99%)	-0.23	14 (1%) 75 70	28, 34, 42, 50	0
1	B	1024/1032 (99%)	-0.18	20 (1%) 65 56	28, 34, 42, 50	0
1	C	1024/1032 (99%)	-0.07	26 (2%) 57 47	28, 34, 42, 50	0
1	D	1024/1032 (99%)	-0.16	17 (1%) 70 63	28, 34, 42, 50	0
All	All	4096/4128 (99%)	-0.16	77 (1%) 66 59	28, 34, 42, 50	0

The worst 5 of 77 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	272	SER	9.5
1	C	270	THR	9.3
1	D	273	THR	9.2
1	D	270	THR	8.5
1	C	269	THR	8.4

### 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
4	NA	D	3006	1/1	0.65	0.22	36,36,36,36	0
2	GAL	B	2001	12/12	0.89	0.17	35,38,39,41	0
4	NA	C	3006	1/1	0.89	0.23	33,33,33,33	0
2	GAL	C	2001	12/12	0.90	0.17	39,45,45,46	0
2	GAL	A	2001	12/12	0.92	0.15	39,41,42,42	0
3	MG	A	3001	1/1	0.92	0.28	22,22,22,22	0
2	GAL	D	2001	12/12	0.93	0.17	41,44,45,45	0
3	MG	C	3001	1/1	0.94	0.16	38,38,38,38	0
4	NA	B	3003	1/1	0.94	0.11	13,13,13,13	0
4	NA	B	3006	1/1	0.95	0.11	19,19,19,19	0
4	NA	A	3006	1/1	0.95	0.07	22,22,22,22	0
3	MG	B	3001	1/1	0.95	0.34	30,30,30,30	0
4	NA	B	3004	1/1	0.96	0.22	5,5,5,5	0
5	MN	D	3005	1/1	0.96	0.06	25,25,25,25	0
4	NA	A	3002	1/1	0.97	0.14	2,2,2,2	0
4	NA	C	3002	1/1	0.97	0.20	11,11,11,11	0
4	NA	D	3002	1/1	0.97	0.23	9,9,9,9	0
4	NA	D	3003	1/1	0.97	0.15	20,20,20,20	0
4	NA	B	3002	1/1	0.98	0.21	2,2,2,2	0
4	NA	A	3003	1/1	0.98	0.11	19,19,19,19	0
4	NA	C	3003	1/1	0.98	0.07	18,18,18,18	0
4	NA	C	3004	1/1	0.98	0.17	13,13,13,13	0
3	MG	D	3001	1/1	0.98	0.12	23,23,23,23	0
4	NA	D	3004	1/1	0.98	0.10	5,5,5,5	0
5	MN	A	3005	1/1	0.99	0.05	24,24,24,24	0
5	MN	B	3005	1/1	0.99	0.07	30,30,30,30	0
4	NA	A	3004	1/1	0.99	0.07	2,2,2,2	0
5	MN	C	3005	1/1	0.99	0.05	30,30,30,30	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.