

wwPDB X-ray Structure Validation Summary Report (i)

Aug 8, 2023 – 08:36 AM EDT

PDB ID : 1NS7

Title: Crystal structure of galactose mutarotase from Lactococcus lactis mutant

E304A complexed with glucose

Authors: Holden, H.M.; Thoden, J.B.

Deposited on : 2003-01-27

Resolution : 1.85 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 (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 (1)) were used in the production of this report:

Mol Probity : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.35

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

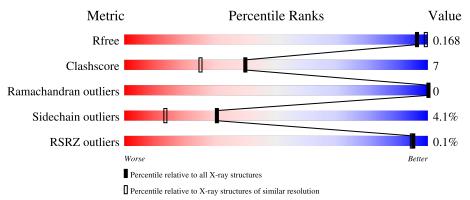
Validation Pipeline (wwPDB-VP) : 2.35

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 1.85 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	Similar resolution
Metric	$(\# ext{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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	347	67%	27%	
1	В	347	73%	22%	•



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5814 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 GALACTOSE MUTAROTASE.

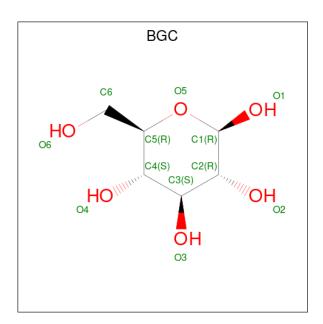
\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	339	Total	С	N	О	S	0	1	0
-	11	330	2645	1669	447	526	3	U	1	
1	D	346	Total	С	N	Ο	S	0	9	0
1	Б	340	2723	1715	466	539	3			U

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	SER	GLU	cloning artifact	UNP Q9ZB17
A	304	ALA	GLU	engineered mutation	UNP Q9ZB17
A	340	LEU	-	expression tag	UNP Q9ZB17
A	341	GLU	-	expression tag	UNP Q9ZB17
A	342	HIS	-	expression tag	UNP Q9ZB17
A	343	HIS	-	expression tag	UNP Q9ZB17
A	344	HIS	-	expression tag	UNP Q9ZB17
A	345	HIS	-	expression tag	UNP Q9ZB17
A	346	HIS	-	expression tag	UNP Q9ZB17
A	347	HIS	-	expression tag	UNP Q9ZB17
В	2	SER	GLU	cloning artifact	UNP Q9ZB17
В	304	ALA	GLU	engineered mutation	UNP Q9ZB17
В	340	LEU	_	expression tag	UNP Q9ZB17
В	341	GLU	-	expression tag	UNP Q9ZB17
В	342	HIS	_	expression tag	UNP Q9ZB17
В	343	HIS	-	expression tag	UNP Q9ZB17
В	344	HIS	-	expression tag	UNP Q9ZB17
В	345	HIS	-	expression tag	UNP Q9ZB17
В	346	HIS	-	expression tag	UNP Q9ZB17
В	347	HIS	_	expression tag	UNP Q9ZB17

• Molecule 2 is beta-D-glucopyranose (three-letter code: BGC) (formula: $C_6H_{12}O_6$).





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

• Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Na 1 1	0	0

• Molecule 4 is water.

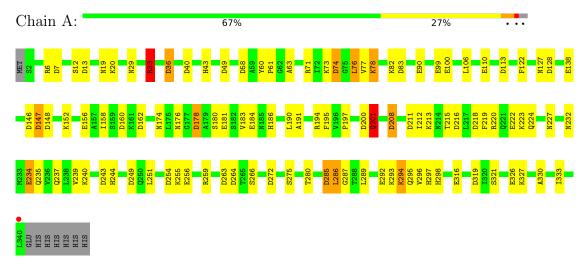
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	197	Total O 197 197	0	0
4	В	224	Total O 224 224	0	0



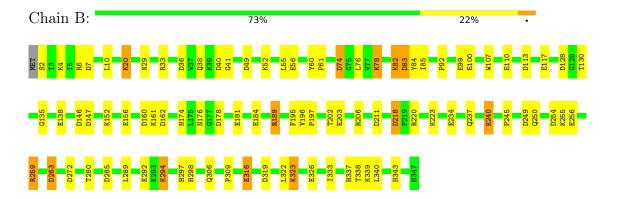
3 Residue-property plots (i)

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: GALACTOSE MUTAROTASE



• Molecule 1: GALACTOSE MUTAROTASE





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	44.90Å 76.20Å 210.90Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 - 1.85	Depositor
resolution (A)	71.67 - 1.81	EDS
% Data completeness	96.3 (30.00-1.85)	Depositor
(in resolution range)	94.1 (71.67-1.81)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) > 1$	2.02 (at 1.82Å)	Xtriage
Refinement program	TNT	Depositor
P.P.	0.167 , 0.228	Depositor
R, R_{free}	0.167 , 0.168	DCC
R_{free} test set	6205 reflections (9.88%)	wwPDB-VP
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.331	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32 , 109.5	EDS
L-test for twinning ²	$ < L >=0.48, < L^2>=0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5814	wwPDB-VP
Average B, all atoms $(Å^2)$	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.71% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: BGC, NA

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	Mol Chain		nd lengths	Bond angles	
WIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5
1	A	0.96	$13/2701 \ (0.5\%)$	1.45	59/3656 (1.6%)
1	В	1.00	$14/2789 \ (0.5\%)$	1.41	44/3774 (1.2%)
All	All	0.98	$27/5490 \ (0.5\%)$	1.43	103/7430 (1.4%)

The worst 5 of 27 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	Ideal(Å)
1	В	234	GLU	CD-OE2	7.19	1.33	1.25
1	A	99	GLU	CD-OE2	6.81	1.33	1.25
1	A	326	GLU	CD-OE2	6.71	1.33	1.25
1	A	100	GLU	CD-OE2	6.40	1.32	1.25
1	В	99	GLU	CD-OE2	6.26	1.32	1.25

The worst 5 of 103 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}(^{o})$
1	В	146	ASP	CB-CG-OD1	10.07	127.36	118.30
1	В	259	ARG	NE-CZ-NH1	9.90	125.25	120.30
1	A	147	ASP	CB-CG-OD2	-9.78	109.49	118.30
1	В	249	ASP	CB-CG-OD2	-9.56	109.69	118.30
1	В	146	ASP	CB-CG-OD2	-9.52	109.73	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	$\mathbf{H}(\mathbf{model})$	H(added)	Clashes	Symm-Clashes
1	A	2645	0	2586	43	0
1	В	2723	0	2637	36	0
2	A	12	0	12	0	0
2	В	12	0	12	0	0
3	A	1	0	0	0	0
4	A	197	0	0	3	0
4	В	224	0	0	4	0
All	All	5814	0	5247	78	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 7.

The worst 5 of 78 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:A:180:SER:HA	1:A:296:VAL:HG13	1.62	0.81
1:A:321:SER:HB2	4:A:1550:HOH:O	1.87	0.74
1:A:20:LYS:HD2	1:A:147:ASP:OD2	1.88	0.73
1:A:232:ASN:HB3	4:A:1562:HOH:O	1.94	0.68
1:B:263:ASP:H	1:B:343:HIS:HD1	1.43	0.66

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	Perce	entiles
1	A	338/347 (97%)	319 (94%)	19 (6%)	0	100	100
1	В	346/347 (100%)	328 (95%)	18 (5%)	0	100	100
All	All	684/694 (99%)	647 (95%)	37 (5%)	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	Percent	iles
1	A	292/299 (98%)	280 (96%)	12 (4%)	30 1	13
1	В	300/299~(100%)	288 (96%)	12 (4%)	31 1	14
All	All	592/598~(99%)	568 (96%)	24 (4%)	30 1	13

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	78	LYS
1	В	195	PHE
1	В	189	ARG
1	В	197	PRO
1	A	224	GLN

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

Mol	Chain	Res	Type
1	A	227	ASN
1	В	38	GLN
1	В	224	GLN
1	В	295	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)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	Trme	Chain	Dag	Link	Bond lengths			Bond angles		
Moi Type	Type		Res		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BGC	A	1400	-	12,12,12	0.65	0	17,17,17	1.13	1 (5%)
2	BGC	В	2400	-	12,12,12	0.37	0	17,17,17	1.37	2 (11%)

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	BGC	A	1400	-	=	0/2/22/22	0/1/1/1
2	BGC	В	2400	-	=	0/2/22/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	В	2400	BGC	C3-C4-C5	-2.95	104.98	110.24
2	В	2400	BGC	O2-C2-C1	2.53	115.03	109.16
2	A	1400	BGC	O2-C2-C1	2.08	113.98	109.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ>2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	$OWAB(A^2)$	Q < 0.9
1	A	339/347 (97%)	-0.75	1 (0%) 94 93	13, 26, 58, 91	0
1	В	346/347 (99%)	-0.81	0 100 100	14, 23, 49, 64	0
All	All	685/694 (98%)	-0.78	1 (0%) 95 95	13, 24, 55, 91	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	340	LEU	2.5

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)

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^{th} 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	${f B-factors}({f \AA}^2)$	Q<0.9
2	BGC	A	1400	12/12	0.91	0.11	22,34,51,64	0
2	BGC	В	2400	12/12	0.94	0.09	19,29,60,68	0
3	NA	A	1401	1/1	0.98	0.11	16,16,16,16	0



6.5 Other polymers (i)

There are no such residues in this entry.

