



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 12:06 AM EDT

PDB ID : 1NS0 / pdb_00001ns0
Title : Crystal structure of galactose mutarotase from *Lactococcus lactis* mutant E304Q complexed with galactose
Authors : Holden, H.M.; Thoden, J.B.
Deposited on : 2003-01-27
Resolution : 1.85 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 ⓘ 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

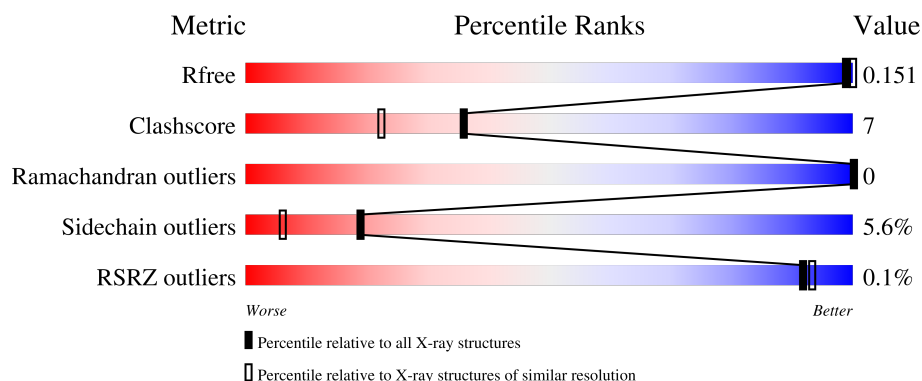
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (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 $\leq 5\%$. 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	347	 77% 18% . .
1	B	347	 79% 17% .

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
2	GLA	A	1400[A]	X	-	-	-
2	GLA	A	1400[B]	X	-	-	-
2	GLA	B	2400[A]	X	-	-	-
2	GLA	B	2400[B]	X	-	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5862 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.

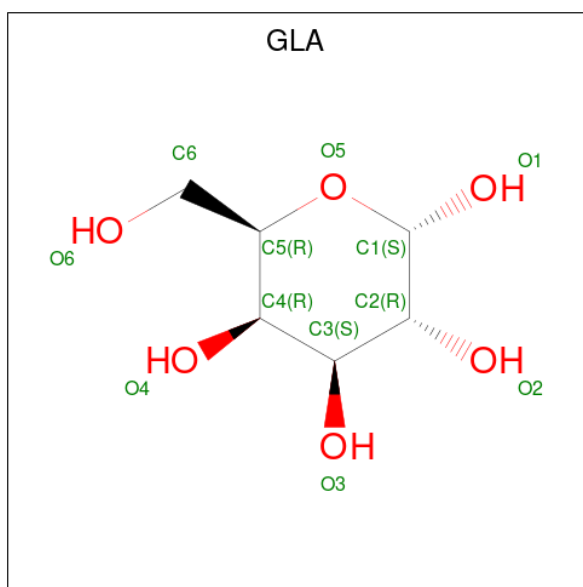
- Molecule 1 is a protein called GALACTOSE MUTAROTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	1	0
			2658	1676	449	530	3			
1	B	346	Total	C	N	O	S	0	1	0
			2722	1714	467	538	3			

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	GLN	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
B	2	SER	GLU	cloning artifact	UNP Q9ZB17
B	304	GLN	GLU	engineered mutation	UNP Q9ZB17
B	340	LEU	-	expression tag	UNP Q9ZB17
B	341	GLU	-	expression tag	UNP Q9ZB17
B	342	HIS	-	expression tag	UNP Q9ZB17
B	343	HIS	-	expression tag	UNP Q9ZB17
B	344	HIS	-	expression tag	UNP Q9ZB17
B	345	HIS	-	expression tag	UNP Q9ZB17
B	346	HIS	-	expression tag	UNP Q9ZB17
B	347	HIS	-	expression tag	UNP Q9ZB17

- Molecule 2 is alpha-D-galactopyranose (CCD ID: GLA) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	1
			13	6	7		
2	B	1	Total	C	O	0	1
			13	6	7		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

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

- Molecule 4 is water.

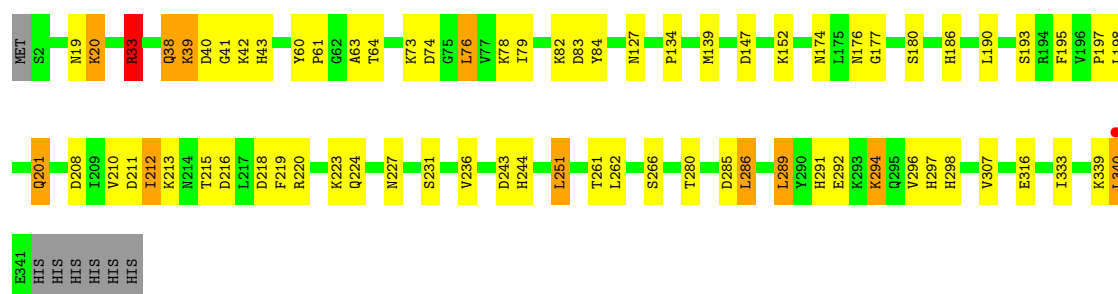
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	211	Total	O	0	0
			211	211		
4	B	244	Total	O	0	0
			244	244		

3 Residue-property plots

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.

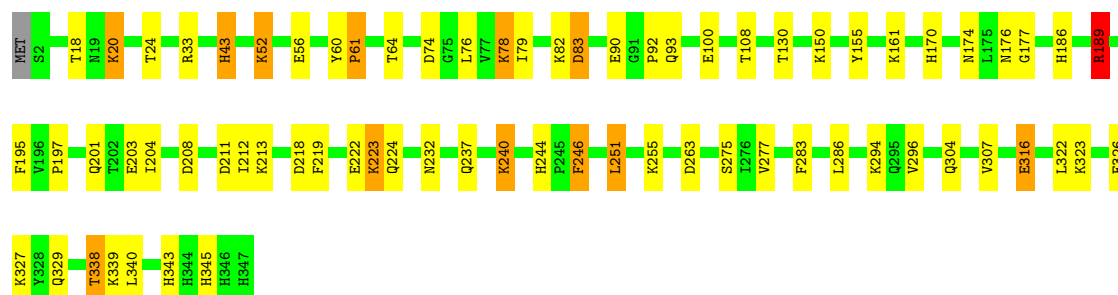
• Molecule 1: GALACTOSE MUTAROTASE

Chain A: 



• Molecule 1: GALACTOSE MUTAROTASE

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	44.80Å 76.20Å 211.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.85 30.00 – 1.85	Depositor EDS
% Data completeness (in resolution range)	96.7 (30.00-1.85) 96.4 (30.00-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 1.81Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.160 , 0.210 0.155 , 0.151	Depositor DCC
R_{free} test set	6299 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å ²)	20.5	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 82.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5862	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 4.63% 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

5.1 Standard geometry

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

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	1.08	0/2714	1.54	17/3673 (0.5%)
1	B	1.08	0/2784	1.50	22/3767 (0.6%)
All	All	1.08	0/5498	1.52	39/7440 (0.5%)

There are no bond length outliers.

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	340	LEU	O-C-N	9.58	134.32	122.63
1	A	340	LEU	CA-C-N	8.76	137.46	121.70
1	A	340	LEU	C-N-CA	8.76	137.46	121.70
1	B	43	HIS	CB-CA-C	-7.36	99.22	110.26
1	A	340	LEU	N-CA-CB	7.28	122.62	110.53
1	A	38	GLN	N-CA-C	7.10	120.24	109.23
1	A	76	LEU	N-CA-C	7.09	120.29	108.73
1	B	244	HIS	O-C-N	7.01	126.19	121.71
1	B	33	ARG	N-CA-CB	-6.96	99.18	111.08
1	B	211	ASP	CA-CB-CG	-6.83	105.77	112.60
1	B	189	ARG	N-CA-CB	6.76	122.03	110.68
1	B	186	HIS	CA-CB-CG	6.68	120.48	113.80
1	A	33	ARG	N-CA-C	6.62	120.31	109.72
1	A	244	HIS	O-C-N	6.42	126.35	121.85
1	A	210	VAL	N-CA-C	6.08	116.63	108.11
1	B	246	PHE	CA-CB-CG	-5.75	108.05	113.80
1	B	203	GLU	N-CA-C	5.74	120.23	113.23
1	A	307	VAL	N-CA-C	-5.74	101.41	109.21
1	B	108	THR	N-CA-C	-5.71	102.85	110.55
1	B	76	LEU	N-CA-C	5.69	118.01	108.73
1	B	100	GLU	CB-CA-C	-5.66	103.85	111.89
1	A	244	HIS	CB-CA-C	-5.63	100.72	109.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	345	HIS	CA-CB-CG	-5.62	108.18	113.80
1	B	232	ASN	N-CA-CB	5.58	119.38	110.73
1	B	307	VAL	N-CA-C	-5.42	101.74	108.89
1	A	33	ARG	N-CA-CB	-5.39	101.86	111.08
1	B	83	ASP	CA-CB-CG	5.36	117.96	112.60
1	B	204	ILE	N-CA-C	-5.35	104.22	110.05
1	A	289	LEU	N-CA-C	5.34	117.44	108.73
1	B	170	HIS	N-CA-C	5.31	119.47	112.25
1	B	33	ARG	N-CA-C	5.29	118.19	109.72
1	A	134	PRO	N-CA-CB	5.25	108.21	103.33
1	A	244	HIS	N-CA-C	5.24	116.72	108.23
1	B	64	THR	N-CA-C	-5.22	100.65	108.96
1	B	296	VAL	CB-CA-C	-5.22	104.28	111.49
1	B	61	PRO	N-CA-CB	5.13	108.63	103.25
1	A	64	THR	N-CA-C	-5.10	101.45	109.25
1	A	236	VAL	CA-C-O	5.07	126.22	120.85
1	B	283	PHE	CA-CB-CG	-5.01	108.79	113.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2658	0	2595	41	0
1	B	2722	0	2638	38	0
2	A	13	0	4	0	0
2	B	13	0	4	0	0
3	A	1	0	0	0	0
4	A	211	0	0	0	0
4	B	244	0	0	5	0
All	All	5862	0	5241	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.

All (78) 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:A:20:LYS:HD2	1:A:147:ASP:OD2	1.92	0.69
1:A:33:ARG:HB2	1:A:63:ALA:HA	1.74	0.67
1:B:189:ARG:HG3	1:B:189:ARG:NH1	2.13	0.64
1:B:189:ARG:HD3	4:B:2578:HOH:O	1.98	0.62
1:B:52:LYS:HE2	1:B:56:GLU:CD	2.24	0.61
1:A:201:GLN:H	1:A:201:GLN:CD	2.09	0.60
1:A:212:ILE:O	1:A:215:THR:HG23	2.02	0.60
1:B:316:GLU:H	1:B:316:GLU:CD	2.09	0.59
1:B:189:ARG:HG3	1:B:189:ARG:HH11	1.68	0.59
1:A:180:SER:HA	1:A:296:VAL:HG13	1.86	0.57
1:A:251:LEU:HD23	1:A:251:LEU:N	2.19	0.57
1:A:174:ASN:OD1	1:A:176:ASN:HB2	2.05	0.57
1:B:263:ASP:H	1:B:343:HIS:HD1	1.53	0.57
1:B:322:LEU:HD12	1:B:326[B]:GLU:HG2	1.87	0.56
1:B:24:THR:HB	4:B:2581:HOH:O	2.05	0.56
1:A:152:LYS:HG3	1:A:333:ILE:HG12	1.87	0.56
1:B:237:GLN:O	1:B:240:LYS:HE2	2.06	0.55
1:A:218:ASP:OD1	1:A:220:ARG:HD3	2.08	0.53
1:A:201:GLN:H	1:A:201:GLN:NE2	2.08	0.51
1:A:38:GLN:HB2	1:A:42:LYS:O	2.11	0.51
1:B:338:THR:HG22	1:B:339:LYS:HG3	1.92	0.50
1:B:322:LEU:CD1	1:B:326[B]:GLU:HG2	2.42	0.50
1:A:219:PHE:CE2	1:A:223:LYS:HG2	2.47	0.49
1:B:74:ASP:HA	1:B:92:PRO:O	2.11	0.49
1:A:19:ASN:HB2	1:A:147:ASP:OD1	2.11	0.49
1:A:291:HIS:O	1:A:292:GLU:HB2	2.13	0.48
1:A:280:THR:O	1:A:298:HIS:HA	2.13	0.48
1:A:297:HIS:O	1:A:298:HIS:HB2	2.14	0.48
1:B:82:LYS:HB3	1:B:82:LYS:HE3	1.62	0.48
1:A:79:ILE:HB	1:A:84:TYR:CE1	2.49	0.47
1:B:174:ASN:OD1	1:B:176:ASN:HB2	2.14	0.47
1:A:197:PRO:HD2	1:A:208:ASP:O	2.15	0.47
1:B:43:HIS:O	1:B:177:GLY:HA2	2.15	0.47
1:B:79:ILE:N	1:B:82:LYS:O	2.43	0.46
1:B:52:LYS:O	1:B:56:GLU:HG3	2.16	0.46
1:A:219:PHE:CD2	1:A:223:LYS:HG2	2.51	0.45
1:A:211:ASP:OD1	1:A:213:LYS:HB2	2.16	0.45
1:B:219:PHE:CD2	1:B:223:LYS:HG2	2.52	0.45
1:A:60:TYR:N	1:A:61:PRO:CD	2.80	0.45
1:B:224:GLN:NE2	4:B:2640:HOH:O	2.46	0.45
1:A:198:LEU:HD11	1:A:243:ASP:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:GLU:HG2	1:B:223:LYS:N	2.31	0.45
1:A:78:LYS:HG2	1:A:83:ASP:OD1	2.17	0.44
1:A:127:ASN:HB2	1:B:130:THR:HG23	1.99	0.44
1:B:60:TYR:N	1:B:61:PRO:CD	2.80	0.44
1:A:39:LYS:O	1:A:40:ASP:C	2.60	0.44
1:A:82:LYS:HD3	1:A:84:TYR:CE1	2.53	0.44
1:A:289:LEU:HD23	1:A:294:LYS:HA	1.98	0.44
1:B:189:ARG:HH11	1:B:189:ARG:CG	2.30	0.44
1:B:161:LYS:HE3	4:B:2575:HOH:O	2.17	0.44
1:A:285:ASP:O	1:A:286:LEU:C	2.60	0.43
1:B:150:LYS:NZ	4:B:2641:HOH:O	2.37	0.43
1:B:78:LYS:HG3	1:B:83:ASP:OD1	2.19	0.43
1:A:20:LYS:HD2	1:A:147:ASP:CG	2.44	0.43
1:B:197:PRO:HD2	1:B:208:ASP:O	2.18	0.43
1:B:277:VAL:HB	1:B:304:GLN:HB2	1.99	0.43
1:A:73:LYS:O	1:A:74:ASP:HB2	2.19	0.43
1:B:218:ASP:O	1:B:223:LYS:HE2	2.18	0.43
1:A:193:SER:HA	1:A:212:ILE:HD11	2.01	0.42
1:B:90:GLU:O	1:B:93:GLN:HB2	2.19	0.42
1:B:155:TYR:O	1:B:329:GLN:HA	2.19	0.42
1:B:316:GLU:CD	1:B:316:GLU:N	2.76	0.42
1:A:216:ASP:OD1	1:A:231:SER:HB2	2.20	0.42
1:A:261:THR:HG22	1:A:262:LEU:N	2.35	0.42
1:B:20:LYS:HE3	1:B:20:LYS:HB2	1.74	0.42
1:B:251:LEU:N	1:B:251:LEU:HD23	2.33	0.41
1:A:190:LEU:HB3	1:A:219:PHE:HB2	2.02	0.41
1:A:218:ASP:OD2	1:A:220:ARG:NH1	2.52	0.41
1:B:246:PHE:O	1:B:275:SER:HB2	2.21	0.41
1:B:340:LEU:HA	1:B:340:LEU:HD12	1.77	0.41
1:A:38:GLN:NE2	1:A:41:GLY:HA2	2.36	0.41
1:A:139:MET:HE3	1:A:139:MET:HB2	1.93	0.41
1:B:18:THR:HG23	1:B:24:THR:HG22	2.02	0.41
1:B:286:LEU:HD12	1:B:286:LEU:HA	1.84	0.41
1:A:43:HIS:O	1:A:177:GLY:HA2	2.21	0.40
1:A:218:ASP:CG	1:A:220:ARG:HH11	2.30	0.40
1:A:294:LYS:HB2	1:A:294:LYS:HE2	1.43	0.40
1:A:227:ASN:HD22	1:A:227:ASN:HA	1.61	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	339/347 (98%)	319 (94%)	20 (6%)	0	100	100
1	B	345/347 (99%)	327 (95%)	18 (5%)	0	100	100
All	All	684/694 (99%)	646 (94%)	38 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	294/300 (98%)	278 (95%)	16 (5%)	20	7
1	B	300/300 (100%)	283 (94%)	17 (6%)	18	6
All	All	594/600 (99%)	561 (94%)	33 (6%)	19	6

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LYS
1	A	33	ARG
1	A	39	LYS
1	A	76	LEU
1	A	186	HIS
1	A	195	PHE
1	A	201	GLN
1	A	212	ILE

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Mol	Chain	Res	Type
1	A	224	GLN
1	A	251	LEU
1	A	266	SER
1	A	286	LEU
1	A	294	LYS
1	A	316	GLU
1	A	339	LYS
1	A	340	LEU
1	B	20	LYS
1	B	52	LYS
1	B	78	LYS
1	B	189	ARG
1	B	195	PHE
1	B	201	GLN
1	B	212	ILE
1	B	213	LYS
1	B	223	LYS
1	B	240	LYS
1	B	251	LEU
1	B	255	LYS
1	B	294	LYS
1	B	316	GLU
1	B	323	LYS
1	B	327	LYS
1	B	338	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	88	GLN
1	A	201	GLN
1	A	224	GLN
1	A	227	ASN
1	A	295	GLN
1	B	38	GLN
1	B	43	HIS
1	B	224	GLN
1	B	227	ASN
1	B	304	GLN
1	B	306	GLN
1	B	313	GLN

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Mol	Chain	Res	Type
1	B	337	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is 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	GLA	A	1400[B]	-	12,12,12	0.53	0	17,17,17	1.47	2 (11%)
2	GLA	A	1400[A]	-	12,12,12	0.53	0	17,17,17	1.39	2 (11%)
2	GLA	B	2400[A]	-	12,12,12	0.45	0	17,17,17	1.27	1 (5%)
2	GLA	B	2400[B]	-	12,12,12	0.50	0	17,17,17	1.32	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	GLA	A	1400[B]	-	1/1/5/5	2/2/22/22	0/1/1/1
2	GLA	A	1400[A]	-	1/1/5/5	2/2/22/22	0/1/1/1
2	GLA	B	2400[A]	-	1/1/5/5	2/2/22/22	0/1/1/1
2	GLA	B	2400[B]	-	1/1/5/5	2/2/22/22	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2400[A]	GLA	O5-C1-C2	-3.53	104.09	110.30
2	B	2400[B]	GLA	O5-C1-C2	-3.53	104.09	110.30
2	A	1400[A]	GLA	O2-C2-C1	3.50	117.32	109.25
2	A	1400[B]	GLA	O2-C2-C1	3.50	117.32	109.25
2	A	1400[B]	GLA	O1-C1-C2	3.43	118.93	108.98
2	A	1400[A]	GLA	O1-C1-C2	2.52	116.29	108.98

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1400[A]	GLA	C1
2	A	1400[B]	GLA	C1
2	B	2400[A]	GLA	C1
2	B	2400[B]	GLA	C1

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1400[A]	GLA	O5-C5-C6-O6
2	A	1400[B]	GLA	O5-C5-C6-O6
2	B	2400[A]	GLA	O5-C5-C6-O6
2	B	2400[B]	GLA	O5-C5-C6-O6
2	A	1400[A]	GLA	C4-C5-C6-O6
2	A	1400[B]	GLA	C4-C5-C6-O6
2	B	2400[A]	GLA	C4-C5-C6-O6
2	B	2400[B]	GLA	C4-C5-C6-O6

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 ⓘ

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	340/347 (97%)	-0.59	1 (0%) 90 92	15, 26, 58, 94	1 (0%)
1	B	346/347 (99%)	-0.77	0 100 100	16, 23, 49, 63	1 (0%)
All	All	686/694 (98%)	-0.68	1 (0%) 92 93	15, 24, 55, 94	2 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	340	LEU	2.3

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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	NA	A	1401	1/1	0.89	0.12	43,43,43,43	0
2	GLA	B	2400[B]	12/12	0.97	0.06	16,22,35,44	1
2	GLA	B	2400[A]	12/12	0.97	0.06	16,22,34,35	1
2	GLA	A	1400[B]	12/12	0.98	0.05	17,23,29,40	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GLA	A	1400[A]	12/12	0.98	0.05	17,23,29,40	1

6.5 Other polymers [i](#)

There are no such residues in this entry.