



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:57 PM GMT

PDB ID : 5BXS
Title : LNBase in complex with LNB-NHAcCAS
Authors : Ito, T.; Arakawa, T.; Fushinobu, S.
Deposited on : 2015-06-09
Resolution : 2.20 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20026982

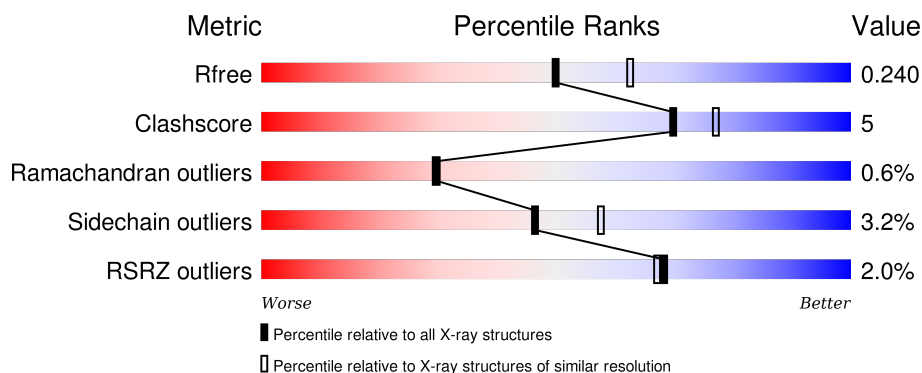
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	644	
1	B	644	

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	GAL	A	701	-	-	-	X
3	GC2	A	702	-	-	-	X
4	SO4	B	703	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10407 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 Lacto-N-biosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	633	Total	C	N	O	S	0	0	0
			4969	3132	846	974	17			
1	B	632	Total	C	N	O	S	0	0	0
			4963	3129	845	972	17			

There are 42 discrepancies between the modelled and reference sequences:

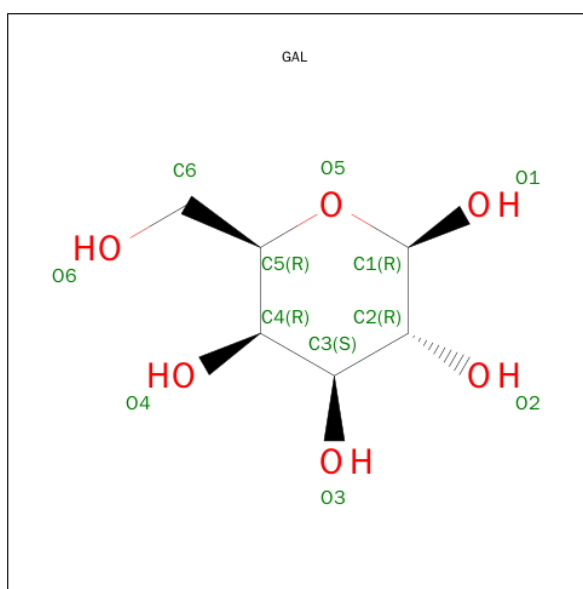
Chain	Residue	Modelled	Actual	Comment	Reference
A	20	MET	-	initiating methionine	UNP B3TLD6
A	21	GLY	-	expression tag	UNP B3TLD6
A	22	SER	-	expression tag	UNP B3TLD6
A	23	SER	-	expression tag	UNP B3TLD6
A	24	HIS	-	expression tag	UNP B3TLD6
A	25	HIS	-	expression tag	UNP B3TLD6
A	26	HIS	-	expression tag	UNP B3TLD6
A	27	HIS	-	expression tag	UNP B3TLD6
A	28	HIS	-	expression tag	UNP B3TLD6
A	29	HIS	-	expression tag	UNP B3TLD6
A	30	SER	-	expression tag	UNP B3TLD6
A	31	SER	-	expression tag	UNP B3TLD6
A	32	GLY	-	expression tag	UNP B3TLD6
A	33	LEU	-	expression tag	UNP B3TLD6
A	34	VAL	-	expression tag	UNP B3TLD6
A	35	PRO	-	expression tag	UNP B3TLD6
A	36	ARG	-	expression tag	UNP B3TLD6
A	37	GLY	-	expression tag	UNP B3TLD6
A	38	SER	-	expression tag	UNP B3TLD6
A	39	HIS	-	expression tag	UNP B3TLD6
A	40	MET	-	expression tag	UNP B3TLD6
B	20	MET	-	initiating methionine	UNP B3TLD6
B	21	GLY	-	expression tag	UNP B3TLD6
B	22	SER	-	expression tag	UNP B3TLD6
B	23	SER	-	expression tag	UNP B3TLD6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	24	HIS	-	expression tag	UNP B3TLD6
B	25	HIS	-	expression tag	UNP B3TLD6
B	26	HIS	-	expression tag	UNP B3TLD6
B	27	HIS	-	expression tag	UNP B3TLD6
B	28	HIS	-	expression tag	UNP B3TLD6
B	29	HIS	-	expression tag	UNP B3TLD6
B	30	SER	-	expression tag	UNP B3TLD6
B	31	SER	-	expression tag	UNP B3TLD6
B	32	GLY	-	expression tag	UNP B3TLD6
B	33	LEU	-	expression tag	UNP B3TLD6
B	34	VAL	-	expression tag	UNP B3TLD6
B	35	PRO	-	expression tag	UNP B3TLD6
B	36	ARG	-	expression tag	UNP B3TLD6
B	37	GLY	-	expression tag	UNP B3TLD6
B	38	SER	-	expression tag	UNP B3TLD6
B	39	HIS	-	expression tag	UNP B3TLD6
B	40	MET	-	expression tag	UNP B3TLD6

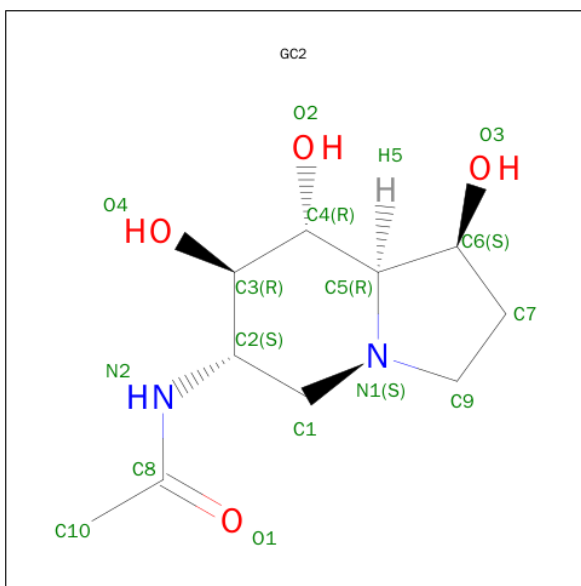
- Molecule 2 is BETA-D-GALACTOSE (three-letter code: GAL) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			11	6	5		
2	B	1	Total	C	O	0	0
			11	6	5		

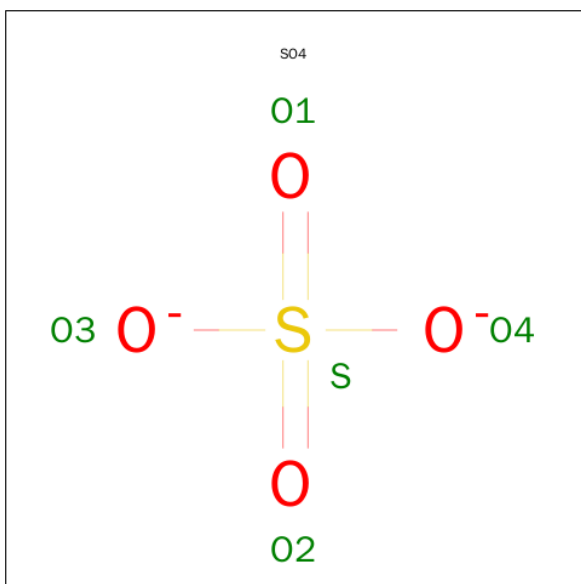
- Molecule 3 is 6-ACETAMIDO-6-DEOXY-CASTANOSPERMINE (three-letter code: GC2)

(formula: C₁₀H₁₈N₂O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			16	10	2	4		
3	B	1	Total	C	N	O	0	0
			16	10	2	4		

- Molecule 4 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		

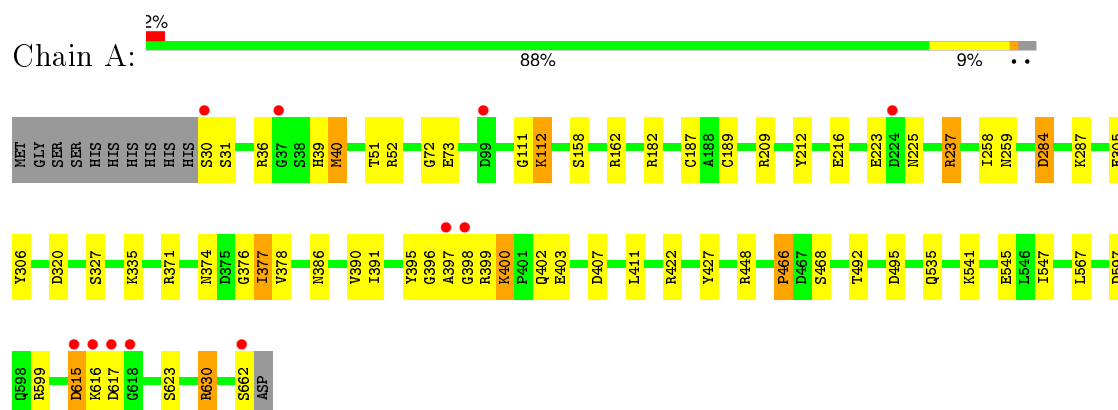
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	200	Total	O	0	0
			200	200		
5	B	211	Total	O	0	0
			211	211		

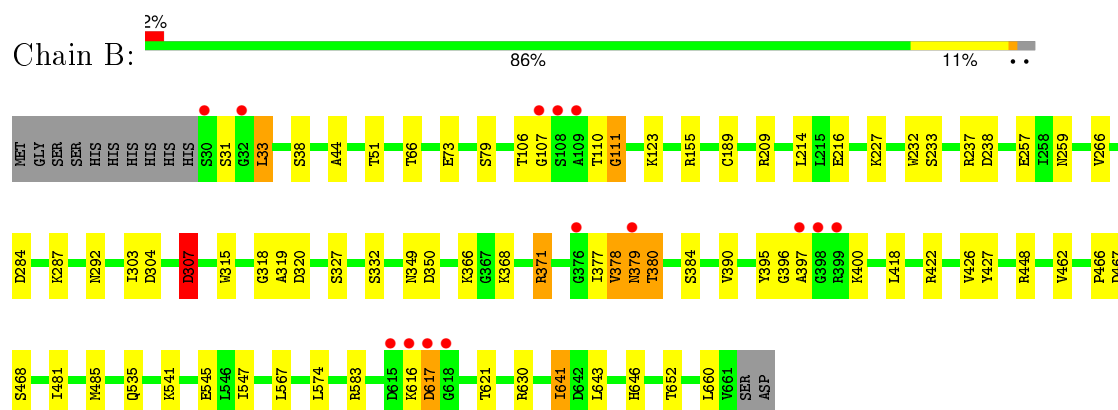
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Lacto-N-biosidase



• Molecule 1: Lacto-N-biosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	116.68Å 131.70Å 104.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.85 – 2.20 38.93 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.85-2.20) 99.7 (38.93-2.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.27 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.191 , 0.238 0.198 , 0.240	Depositor DCC
R_{free} test set	4119 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	19.9	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 38.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 82413 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10407	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GC2, GAL, SO4

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	0.87	1/5078 (0.0%)	0.91	10/6891 (0.1%)
1	B	0.86	0/5072	0.90	5/6883 (0.1%)
All	All	0.86	1/10150 (0.0%)	0.91	15/13774 (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	3
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	597	ASP	CB-CG	6.44	1.65	1.51

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	237	ARG	NE-CZ-NH2	-6.95	116.82	120.30
1	A	52	ARG	NE-CZ-NH1	6.50	123.55	120.30
1	B	237	ARG	NE-CZ-NH2	6.41	123.50	120.30
1	A	284	ASP	CB-CG-OD1	5.51	123.25	118.30
1	B	448	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	307	ASP	CB-CG-OD2	-5.43	113.41	118.30
1	A	599	ARG	NE-CZ-NH1	5.35	122.97	120.30
1	A	225	ASN	N-CA-CB	-5.26	101.12	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	237	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	422	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	B	155	ARG	NE-CZ-NH2	5.21	122.91	120.30
1	B	371	ARG	NE-CZ-NH1	-5.19	117.71	120.30
1	A	630	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	A	237	ARG	CG-CD-NE	-5.15	100.98	111.80
1	A	599	ARG	NE-CZ-NH2	-5.10	117.75	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	468	SER	Peptide
1	B	378	VAL	Peptide
1	B	379	ASN	Peptide
1	B	468	SER	Peptide

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	4969	0	4843	42	0
1	B	4963	0	4838	51	0
2	A	11	0	10	1	0
2	B	11	0	10	1	0
3	A	16	0	17	2	0
3	B	16	0	17	2	0
4	A	5	0	0	0	0
4	B	5	0	0	1	0
5	A	200	0	0	4	0
5	B	211	0	0	2	0
All	All	10407	0	9735	89	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 5.

All (89) 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:B:379:ASN:HB3	1:B:380:THR:HA	1.15	1.13
1:B:378:VAL:HG23	1:B:379:ASN:HB2	1.25	1.09
1:B:379:ASN:HB3	1:B:380:THR:CA	1.82	1.07
1:A:40:MET:HG3	1:A:112:LYS:HE3	1.48	0.95
1:B:378:VAL:HG23	1:B:379:ASN:CB	2.02	0.90
1:A:40:MET:HG3	1:A:112:LYS:CE	2.12	0.79
1:B:238:ASP:HB3	5:B:836:HOH:O	1.83	0.78
1:A:395:TYR:CE2	1:A:397:ALA:HB3	2.23	0.72
1:A:395:TYR:CE2	1:A:397:ALA:CB	2.74	0.71
1:A:427:TYR:CE2	3:A:702:GC2:H72C	2.26	0.70
1:B:189:CYS:HB2	5:B:969:HOH:O	1.91	0.69
1:B:33:LEU:HD23	1:B:33:LEU:C	2.13	0.69
1:B:379:ASN:CB	1:B:380:THR:HA	2.10	0.68
1:A:395:TYR:CD2	1:A:397:ALA:HB3	2.29	0.68
1:B:427:TYR:CE2	3:B:702:GC2:H72C	2.30	0.67
1:A:182:ARG:HG2	1:A:492:THR:HB	1.79	0.65
1:B:259:ASN:ND2	1:B:320:ASP:OD1	2.29	0.62
1:A:376:GLY:O	1:A:378:VAL:HG13	2.00	0.60
1:B:349:ASN:HD22	1:B:379:ASN:ND2	2.00	0.60
1:B:378:VAL:CG2	1:B:379:ASN:HB2	2.17	0.60
1:A:237:ARG:NH2	1:A:305:GLU:O	2.35	0.59
1:A:391:ILE:HB	1:A:411:LEU:HD23	1.86	0.58
1:B:377:ILE:N	1:B:377:ILE:HD13	2.19	0.56
1:B:318:GLY:O	1:B:319:ALA:HB3	2.05	0.56
1:B:379:ASN:CB	1:B:380:THR:CA	2.70	0.56
1:A:623:SER:OG	1:A:630:ARG:HD3	2.07	0.54
1:B:350:ASP:OD1	1:B:380:THR:HB	2.08	0.53
1:B:397:ALA:CA	1:B:400:LYS:HZ3	2.23	0.52
1:A:73:GLU:O	1:A:112:LYS:HE3	2.10	0.52
1:A:187:CYS:HB3	1:A:466:PRO:HD3	1.91	0.52
1:A:189:CYS:HB2	5:A:970:HOH:O	2.09	0.52
1:A:39:HIS:HB3	5:A:989:HOH:O	2.10	0.51
1:B:583:ARG:HD3	1:B:652:THR:HG21	1.94	0.50
1:A:40:MET:CG	1:A:112:LYS:HD3	2.42	0.49
1:A:40:MET:HB3	1:A:72:GLY:HA3	1.93	0.49
1:B:304:ASP:O	1:B:307:ASP:HB2	2.13	0.48
1:A:547:ILE:HG12	5:A:986:HOH:O	2.14	0.48
1:A:31:SER:HA	1:B:535:GLN:HB2	1.95	0.48
1:A:259:ASN:ND2	1:A:320:ASP:OD1	2.41	0.48
1:A:427:TYR:CE2	3:A:702:GC2:C7	2.97	0.47
1:B:621:THR:HB	1:B:630:ARG:HD2	1.97	0.47
1:B:307:ASP:OD2	1:B:366:LYS:HD3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:303:ILE:HG12	1:B:315:TRP:CZ2	2.50	0.47
1:B:616:LYS:O	1:B:617:ASP:CB	2.62	0.47
1:B:371:ARG:HG2	1:B:390:VAL:HB	1.96	0.46
1:A:212:TYR:C	1:A:212:TYR:CD1	2.88	0.46
1:B:33:LEU:C	1:B:33:LEU:CD2	2.83	0.46
1:A:374:ASN:O	1:A:377:ILE:HG22	2.16	0.46
1:B:418:LEU:HB3	1:B:462:VAL:HA	1.97	0.46
1:A:223:GLU:HG3	1:A:305:GLU:OE2	2.16	0.45
1:B:284:ASP:OD2	1:B:287:LYS:NZ	2.45	0.45
1:B:44:ALA:O	1:B:66:THR:HG22	2.17	0.45
1:B:110:THR:HG22	1:B:111:GLY:N	2.32	0.45
1:B:422:ARG:HG3	1:B:466:PRO:HA	1.99	0.45
1:B:349:ASN:ND2	1:B:378:VAL:HG21	2.32	0.44
1:B:616:LYS:O	1:B:617:ASP:HB2	2.17	0.44
1:A:615:ASP:O	1:A:617:ASP:HA	2.17	0.44
1:B:216:GLU:OE2	2:B:701:GAL:O3	2.25	0.44
1:B:643:LEU:HA	1:B:646:HIS:O	2.18	0.43
1:A:40:MET:CG	1:A:112:LYS:CE	2.91	0.43
1:A:216:GLU:OE2	2:A:701:GAL:H3	2.18	0.43
1:A:400:LYS:HE3	1:A:448:ARG:HH21	1.83	0.43
1:A:40:MET:HB2	1:A:112:LYS:CD	2.48	0.43
1:B:427:TYR:CD2	3:B:702:GC2:H72C	2.53	0.43
1:A:399:ARG:HB3	1:A:403:GLU:OE1	2.18	0.43
1:B:378:VAL:HG23	1:B:379:ASN:N	2.33	0.43
1:A:395:TYR:CE2	1:A:397:ALA:HB2	2.52	0.43
1:A:386:ASN:HB3	5:A:976:HOH:O	2.19	0.42
1:B:79:SER:O	1:B:106:THR:HA	2.19	0.42
1:B:467:ASP:O	1:B:574:LEU:HB3	2.20	0.42
1:B:641:ILE:HG13	1:B:641:ILE:O	2.19	0.42
1:B:214:LEU:HD11	1:B:257:GLU:HB2	2.02	0.42
1:A:495:ASP:O	1:A:495:ASP:CG	2.57	0.42
1:A:535:GLN:HB2	1:B:31:SER:HA	2.02	0.41
1:A:545:GLU:OE1	1:B:51:THR:OG1	2.25	0.41
1:B:397:ALA:HA	1:B:400:LYS:HZ3	1.85	0.41
1:A:400:LYS:CE	1:A:448:ARG:HH21	2.33	0.41
1:A:284:ASP:OD2	1:A:287:LYS:NZ	2.48	0.41
1:A:51:THR:OG1	1:B:545:GLU:OE1	2.36	0.41
1:A:374:ASN:O	1:A:377:ILE:CG2	2.69	0.41
1:B:481:ILE:O	1:B:485:MET:HG2	2.20	0.41
1:B:426:VAL:HG13	4:B:703:SO4:O4	2.21	0.41
1:B:292:ASN:OD1	1:B:292:ASN:C	2.59	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:395:TYR:CE2	1:B:397:ALA:HB2	2.56	0.40
1:A:158:SER:O	1:A:162:ARG:HD3	2.21	0.40
1:B:232:TRP:CG	1:B:233:SER:HA	2.55	0.40
1:A:371:ARG:HG2	1:A:390:VAL:HB	2.04	0.40
1:B:397:ALA:HA	1:B:400:LYS:NZ	2.37	0.40
1:A:258:ILE:HD11	1:A:306:TYR:CE1	2.56	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	631/644 (98%)	606 (96%)	21 (3%)	4 (1%)	30	29
1	B	630/644 (98%)	599 (95%)	27 (4%)	4 (1%)	30	29
All	All	1261/1288 (98%)	1205 (96%)	48 (4%)	8 (1%)	30	29

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	111	GLY
1	A	396	GLY
1	B	107	GLY
1	B	396	GLY
1	B	617	ASP
1	A	398	GLY
1	A	111	GLY
1	A	466	PRO

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	530/540 (98%)	514 (97%)	16 (3%)	48	60
1	B	529/540 (98%)	511 (97%)	18 (3%)	44	54
All	All	1059/1080 (98%)	1025 (97%)	34 (3%)	46	57

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	SER
1	A	36	ARG
1	A	40	MET
1	A	112	LYS
1	A	209	ARG
1	A	327	SER
1	A	335	LYS
1	A	377	ILE
1	A	400	LYS
1	A	402	GLN
1	A	407	ASP
1	A	541	LYS
1	A	567	LEU
1	A	615	ASP
1	A	616	LYS
1	A	662	SER
1	B	33	LEU
1	B	38	SER
1	B	73	GLU
1	B	123	LYS
1	B	209	ARG
1	B	227	LYS
1	B	266	VAL
1	B	307	ASP
1	B	327	SER
1	B	332	SER
1	B	368	LYS

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Mol	Chain	Res	Type
1	B	380	THR
1	B	384	SER
1	B	541	LYS
1	B	547	ILE
1	B	567	LEU
1	B	641	ILE
1	B	660	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	349	ASN
1	B	437	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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	A	701	3	11,11,12	0.89	0	15,15,17	1.06	1 (6%)
3	GC2	A	702	2	17,17,17	1.06	0	15,25,25	1.57	4 (26%)
4	SO4	A	703	-	4,4,4	0.61	0	6,6,6	0.76	0
2	GAL	B	701	3	11,11,12	0.97	0	15,15,17	1.76	6 (40%)
3	GC2	B	702	2	17,17,17	1.74	4 (23%)	15,25,25	1.50	1 (6%)
4	SO4	B	703	-	4,4,4	1.03	1 (25%)	6,6,6	0.35	0

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	A	701	3	-	0/2/19/22	0/1/1/1
3	GC2	A	702	2	-	0/4/33/33	0/2/2/2
4	SO4	A	703	-	-	0/0/0/0	0/0/0/0
2	GAL	B	701	3	-	0/2/19/22	0/1/1/1
3	GC2	B	702	2	-	0/4/33/33	0/2/2/2
4	SO4	B	703	-	-	0/0/0/0	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	702	GC2	C6-C5	-4.28	1.49	1.54
3	B	702	GC2	C4-C5	-2.97	1.48	1.53
3	B	702	GC2	O3-C6	-2.19	1.38	1.43
4	B	703	SO4	O4-S	-2.01	1.40	1.47
3	B	702	GC2	C2-N2	2.02	1.49	1.46

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	702	GC2	O3-C6-C7	-4.83	100.81	111.43
3	A	702	GC2	O3-C6-C7	-2.46	106.02	111.43
2	B	701	GAL	O3-C3-C2	-2.34	105.72	110.01
2	A	701	GAL	C6-C5-C4	2.02	118.06	112.99
3	A	702	GC2	O3-C6-C5	2.12	115.64	111.00
2	B	701	GAL	O5-C5-C4	2.14	113.68	110.13
2	B	701	GAL	O4-C4-C5	2.28	115.25	109.23
2	B	701	GAL	O3-C3-C4	2.38	115.72	110.36
3	A	702	GC2	C9-C7-C6	2.60	107.70	103.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	GAL	C1-C2-C3	2.66	112.77	109.55
2	B	701	GAL	C6-C5-C4	2.87	120.19	112.99
3	A	702	GC2	O2-C4-C3	2.98	117.07	110.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	701	GAL	1	0
3	A	702	GC2	2	0
2	B	701	GAL	1	0
3	B	702	GC2	2	0
4	B	703	SO4	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 ⓘ

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	633/644 (98%)	-0.34	11 (1%) 73 72	10, 19, 41, 73	0
1	B	632/644 (98%)	-0.32	14 (2%) 65 64	10, 19, 40, 67	0
All	All	1265/1288 (98%)	-0.33	25 (1%) 68 67	10, 19, 40, 73	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	616	LYS	5.5
1	A	617	ASP	5.1
1	B	107	GLY	5.1
1	A	618	GLY	4.8
1	A	616	LYS	4.4
1	B	617	ASP	4.3
1	B	109	ALA	4.2
1	A	397	ALA	4.2
1	A	398	GLY	4.1
1	B	32	GLY	3.3
1	B	615	ASP	3.2
1	B	108	SER	3.0
1	A	37	GLY	2.9
1	A	99	ASP	2.9
1	A	30	SER	2.8
1	B	379	ASN	2.6
1	A	224	ASP	2.4
1	B	399	ARG	2.3
1	A	615	ASP	2.3
1	B	398	GLY	2.3
1	B	618	GLY	2.2
1	B	30	SER	2.2
1	A	662	SER	2.2
1	B	397	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	376	GLY	2.1

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	GAL	A	701	11/12	0.84	0.15	3.30	26,28,31,31	0
3	GC2	A	702	16/16	0.88	0.14	2.49	21,25,28,29	0
4	SO4	B	703	5/5	0.99	0.11	2.09	19,21,24,25	0
3	GC2	B	702	16/16	0.90	0.14	1.83	23,25,27,27	0
4	SO4	A	703	5/5	0.99	0.10	0.94	24,24,27,28	0
2	GAL	B	701	11/12	0.92	0.11	-0.33	20,22,27,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.