



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 11:08 AM UTC

PDB ID : 3VDA / pdb_00003vda
Title : E. coli (lacZ) beta-galactosidase (N460T)
Authors : Wheatley, R.W.; Kappelhoff, J.C.; Hahn, J.N.; Dugdale, M.L.; Dutkoski, M.J.;
Tamman, S.D.; Fraser, M.E.; Huber, R.E.
Deposited on : 2012-01-04
Resolution : 2.50 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

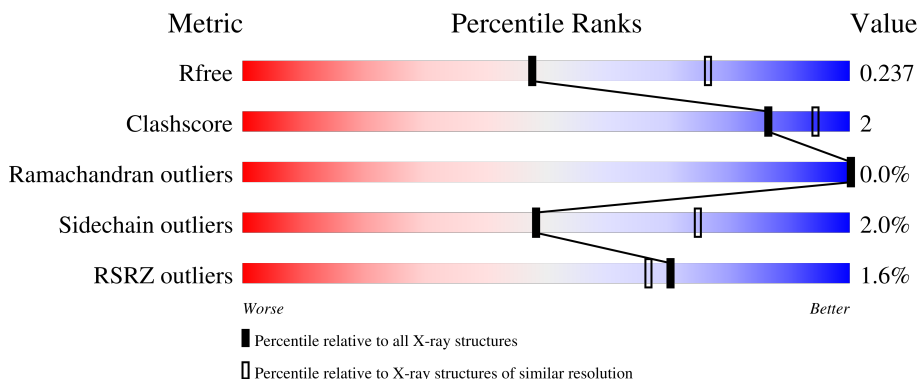
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	1052	% 90% 6% .
1	B	1052	2% 90% 6% .
1	C	1052	2% 89% 7% .
1	D	1052	% 88% 8% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 34381 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
1	A	1015	Total	C	N	O	S	0	0	0
			8156	5159	1444	1515	38			
1	B	1015	Total	C	N	O	S	0	0	0
			8156	5159	1444	1515	38			
1	C	1015	Total	C	N	O	S	0	0	0
			8156	5159	1444	1515	38			
1	D	1015	Total	C	N	O	S	0	0	0
			8156	5159	1444	1515	38			

There are 152 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-28	MET	-	expression tag	UNP P00722
A	-27	GLY	-	expression tag	UNP P00722
A	-26	GLY	-	expression tag	UNP P00722
A	-25	SER	-	expression tag	UNP P00722
A	-24	HIS	-	expression tag	UNP P00722
A	-23	HIS	-	expression tag	UNP P00722
A	-22	HIS	-	expression tag	UNP P00722
A	-21	HIS	-	expression tag	UNP P00722
A	-20	HIS	-	expression tag	UNP P00722
A	-19	HIS	-	expression tag	UNP P00722
A	-18	GLY	-	expression tag	UNP P00722
A	-17	MET	-	expression tag	UNP P00722
A	-16	ALA	-	expression tag	UNP P00722
A	-15	SER	-	expression tag	UNP P00722
A	-14	MET	-	expression tag	UNP P00722
A	-13	THR	-	expression tag	UNP P00722
A	-12	GLY	-	expression tag	UNP P00722
A	-11	GLY	-	expression tag	UNP P00722
A	-10	GLN	-	expression tag	UNP P00722
A	-9	GLN	-	expression tag	UNP P00722
A	-8	MET	-	expression tag	UNP P00722

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	expression tag	UNP P00722
A	-6	ARG	-	expression tag	UNP P00722
A	-5	ASP	-	expression tag	UNP P00722
A	-4	LEU	-	expression tag	UNP P00722
A	-3	TYR	-	expression tag	UNP P00722
A	-2	ASP	-	expression tag	UNP P00722
A	-1	ASP	-	expression tag	UNP P00722
A	0	ASP	-	expression tag	UNP P00722
A	1	ASP	-	expression tag	UNP P00722
A	2	LYS	-	expression tag	UNP P00722
A	3	ASP	-	expression tag	UNP P00722
A	4	PRO	-	expression tag	UNP P00722
A	5	MET	-	expression tag	UNP P00722
A	6	ILE	-	expression tag	UNP P00722
A	7	ASP	-	expression tag	UNP P00722
A	8	PRO	-	expression tag	UNP P00722
A	460	THR	ASN	engineered mutation	UNP P00722
B	-28	MET	-	expression tag	UNP P00722
B	-27	GLY	-	expression tag	UNP P00722
B	-26	GLY	-	expression tag	UNP P00722
B	-25	SER	-	expression tag	UNP P00722
B	-24	HIS	-	expression tag	UNP P00722
B	-23	HIS	-	expression tag	UNP P00722
B	-22	HIS	-	expression tag	UNP P00722
B	-21	HIS	-	expression tag	UNP P00722
B	-20	HIS	-	expression tag	UNP P00722
B	-19	HIS	-	expression tag	UNP P00722
B	-18	GLY	-	expression tag	UNP P00722
B	-17	MET	-	expression tag	UNP P00722
B	-16	ALA	-	expression tag	UNP P00722
B	-15	SER	-	expression tag	UNP P00722
B	-14	MET	-	expression tag	UNP P00722
B	-13	THR	-	expression tag	UNP P00722
B	-12	GLY	-	expression tag	UNP P00722
B	-11	GLY	-	expression tag	UNP P00722
B	-10	GLN	-	expression tag	UNP P00722
B	-9	GLN	-	expression tag	UNP P00722
B	-8	MET	-	expression tag	UNP P00722
B	-7	GLY	-	expression tag	UNP P00722
B	-6	ARG	-	expression tag	UNP P00722
B	-5	ASP	-	expression tag	UNP P00722
B	-4	LEU	-	expression tag	UNP P00722

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	TYR	-	expression tag	UNP P00722
B	-2	ASP	-	expression tag	UNP P00722
B	-1	ASP	-	expression tag	UNP P00722
B	0	ASP	-	expression tag	UNP P00722
B	1	ASP	-	expression tag	UNP P00722
B	2	LYS	-	expression tag	UNP P00722
B	3	ASP	-	expression tag	UNP P00722
B	4	PRO	-	expression tag	UNP P00722
B	5	MET	-	expression tag	UNP P00722
B	6	ILE	-	expression tag	UNP P00722
B	7	ASP	-	expression tag	UNP P00722
B	8	PRO	-	expression tag	UNP P00722
B	460	THR	ASN	engineered mutation	UNP P00722
C	-28	MET	-	expression tag	UNP P00722
C	-27	GLY	-	expression tag	UNP P00722
C	-26	GLY	-	expression tag	UNP P00722
C	-25	SER	-	expression tag	UNP P00722
C	-24	HIS	-	expression tag	UNP P00722
C	-23	HIS	-	expression tag	UNP P00722
C	-22	HIS	-	expression tag	UNP P00722
C	-21	HIS	-	expression tag	UNP P00722
C	-20	HIS	-	expression tag	UNP P00722
C	-19	HIS	-	expression tag	UNP P00722
C	-18	GLY	-	expression tag	UNP P00722
C	-17	MET	-	expression tag	UNP P00722
C	-16	ALA	-	expression tag	UNP P00722
C	-15	SER	-	expression tag	UNP P00722
C	-14	MET	-	expression tag	UNP P00722
C	-13	THR	-	expression tag	UNP P00722
C	-12	GLY	-	expression tag	UNP P00722
C	-11	GLY	-	expression tag	UNP P00722
C	-10	GLN	-	expression tag	UNP P00722
C	-9	GLN	-	expression tag	UNP P00722
C	-8	MET	-	expression tag	UNP P00722
C	-7	GLY	-	expression tag	UNP P00722
C	-6	ARG	-	expression tag	UNP P00722
C	-5	ASP	-	expression tag	UNP P00722
C	-4	LEU	-	expression tag	UNP P00722
C	-3	TYR	-	expression tag	UNP P00722
C	-2	ASP	-	expression tag	UNP P00722
C	-1	ASP	-	expression tag	UNP P00722
C	0	ASP	-	expression tag	UNP P00722

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1	ASP	-	expression tag	UNP P00722
C	2	LYS	-	expression tag	UNP P00722
C	3	ASP	-	expression tag	UNP P00722
C	4	PRO	-	expression tag	UNP P00722
C	5	MET	-	expression tag	UNP P00722
C	6	ILE	-	expression tag	UNP P00722
C	7	ASP	-	expression tag	UNP P00722
C	8	PRO	-	expression tag	UNP P00722
C	460	THR	ASN	engineered mutation	UNP P00722
D	-28	MET	-	expression tag	UNP P00722
D	-27	GLY	-	expression tag	UNP P00722
D	-26	GLY	-	expression tag	UNP P00722
D	-25	SER	-	expression tag	UNP P00722
D	-24	HIS	-	expression tag	UNP P00722
D	-23	HIS	-	expression tag	UNP P00722
D	-22	HIS	-	expression tag	UNP P00722
D	-21	HIS	-	expression tag	UNP P00722
D	-20	HIS	-	expression tag	UNP P00722
D	-19	HIS	-	expression tag	UNP P00722
D	-18	GLY	-	expression tag	UNP P00722
D	-17	MET	-	expression tag	UNP P00722
D	-16	ALA	-	expression tag	UNP P00722
D	-15	SER	-	expression tag	UNP P00722
D	-14	MET	-	expression tag	UNP P00722
D	-13	THR	-	expression tag	UNP P00722
D	-12	GLY	-	expression tag	UNP P00722
D	-11	GLY	-	expression tag	UNP P00722
D	-10	GLN	-	expression tag	UNP P00722
D	-9	GLN	-	expression tag	UNP P00722
D	-8	MET	-	expression tag	UNP P00722
D	-7	GLY	-	expression tag	UNP P00722
D	-6	ARG	-	expression tag	UNP P00722
D	-5	ASP	-	expression tag	UNP P00722
D	-4	LEU	-	expression tag	UNP P00722
D	-3	TYR	-	expression tag	UNP P00722
D	-2	ASP	-	expression tag	UNP P00722
D	-1	ASP	-	expression tag	UNP P00722
D	0	ASP	-	expression tag	UNP P00722
D	1	ASP	-	expression tag	UNP P00722
D	2	LYS	-	expression tag	UNP P00722
D	3	ASP	-	expression tag	UNP P00722
D	4	PRO	-	expression tag	UNP P00722

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Chain	Residue	Modelled	Actual	Comment	Reference
D	5	MET	-	expression tag	UNP P00722
D	6	ILE	-	expression tag	UNP P00722
D	7	ASP	-	expression tag	UNP P00722
D	8	PRO	-	expression tag	UNP P00722
D	460	THR	ASN	engineered mutation	UNP P00722

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		
2	B	2	Total	Mg	0	0
			2	2		
2	C	2	Total	Mg	0	0
			2	2		
2	D	2	Total	Mg	0	0
			2	2		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Na	0	0
			3	3		
3	B	2	Total	Na	0	0
			2	2		
3	C	2	Total	Na	0	0
			2	2		
3	D	3	Total	Na	0	0
			3	3		

- Molecule 4 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			4	2	1	1		
4	A	1	Total	C	O	S	0	0
			4	2	1	1		
4	A	1	Total	C	O	S	0	0
			4	2	1	1		
4	A	1	Total	C	O	S	0	0
			4	2	1	1		
4	A	1	Total	C	O	S	0	0
			4	2	1	1		
4	A	1	Total	C	O	S	0	0
			4	2	1	1		
4	B	1	Total	C	O	S	0	0
			4	2	1	1		
4	B	1	Total	C	O	S	0	0
			4	2	1	1		
4	B	1	Total	C	O	S	0	0
			4	2	1	1		
4	B	1	Total	C	O	S	0	0
			4	2	1	1		
4	B	1	Total	C	O	S	0	0
			4	2	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	O	S	0	0
			4	2	1	1		
4	C	1	Total	C	O	S	0	0
			4	2	1	1		
4	C	1	Total	C	O	S	0	0
			4	2	1	1		
4	C	1	Total	C	O	S	0	0
			4	2	1	1		
4	C	1	Total	C	O	S	0	0
			4	2	1	1		
4	C	1	Total	C	O	S	0	0
			4	2	1	1		
4	C	1	Total	C	O	S	0	0
			4	2	1	1		
4	D	1	Total	C	O	S	0	0
			4	2	1	1		
4	D	1	Total	C	O	S	0	0
			4	2	1	1		
4	D	1	Total	C	O	S	0	0
			4	2	1	1		
4	D	1	Total	C	O	S	0	0
			4	2	1	1		
4	D	1	Total	C	O	S	0	0
			4	2	1	1		
4	D	1	Total	C	O	S	0	0
			4	2	1	1		
4	D	1	Total	C	O	S	0	0
			4	2	1	1		
4	D	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	427	Total	O	0	0
			427	427		

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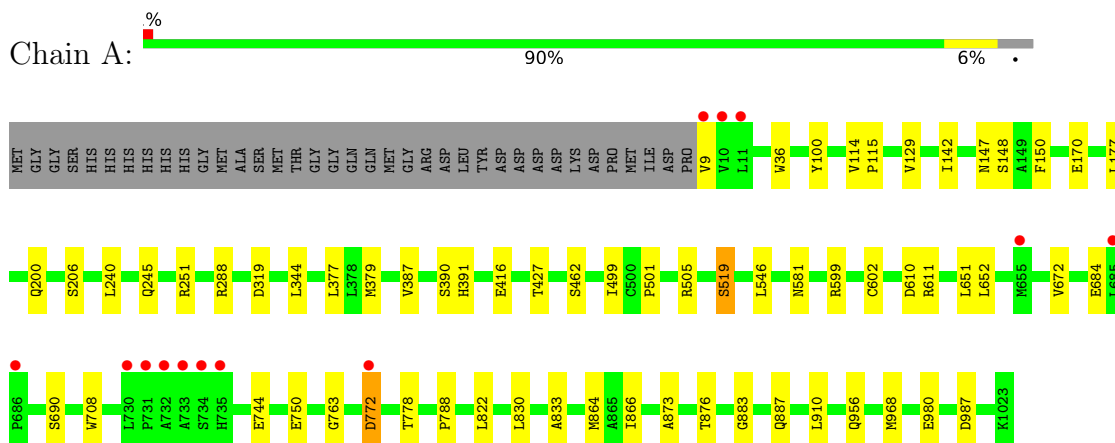
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	403	Total 403	O 403	0	0
5	C	389	Total 389	O 389	0	0
5	D	392	Total 392	O 392	0	0

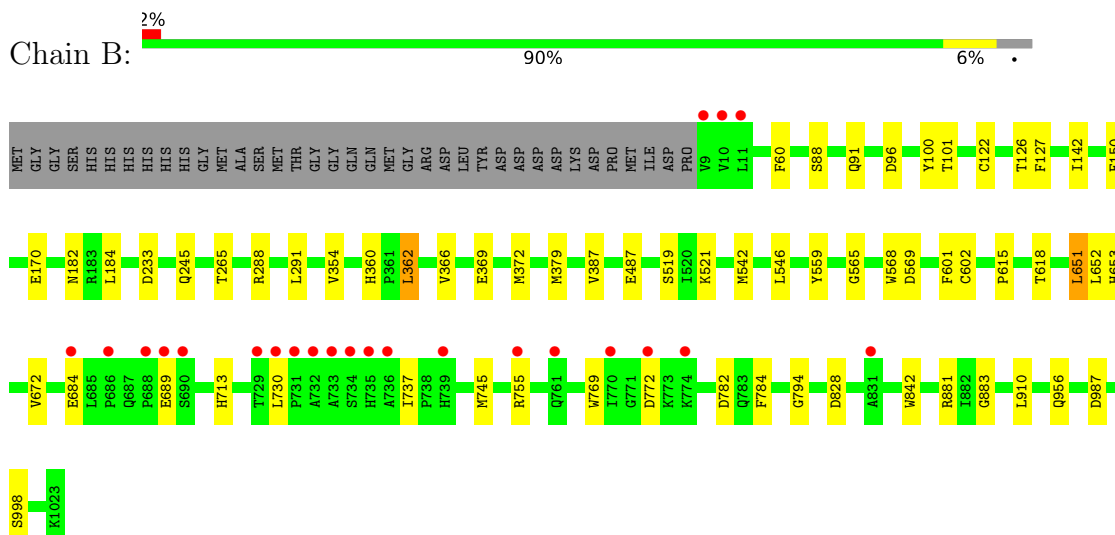
3 Residue-property plots [i](#)

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.

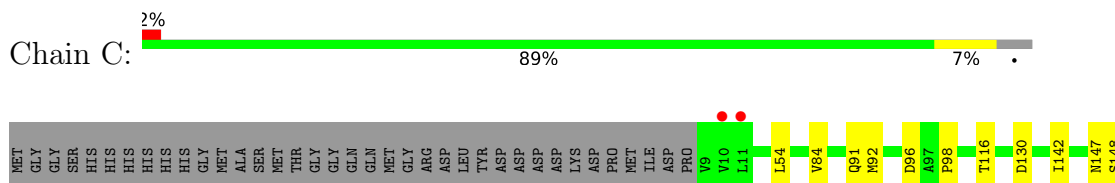
• Molecule 1: Beta-galactosidase

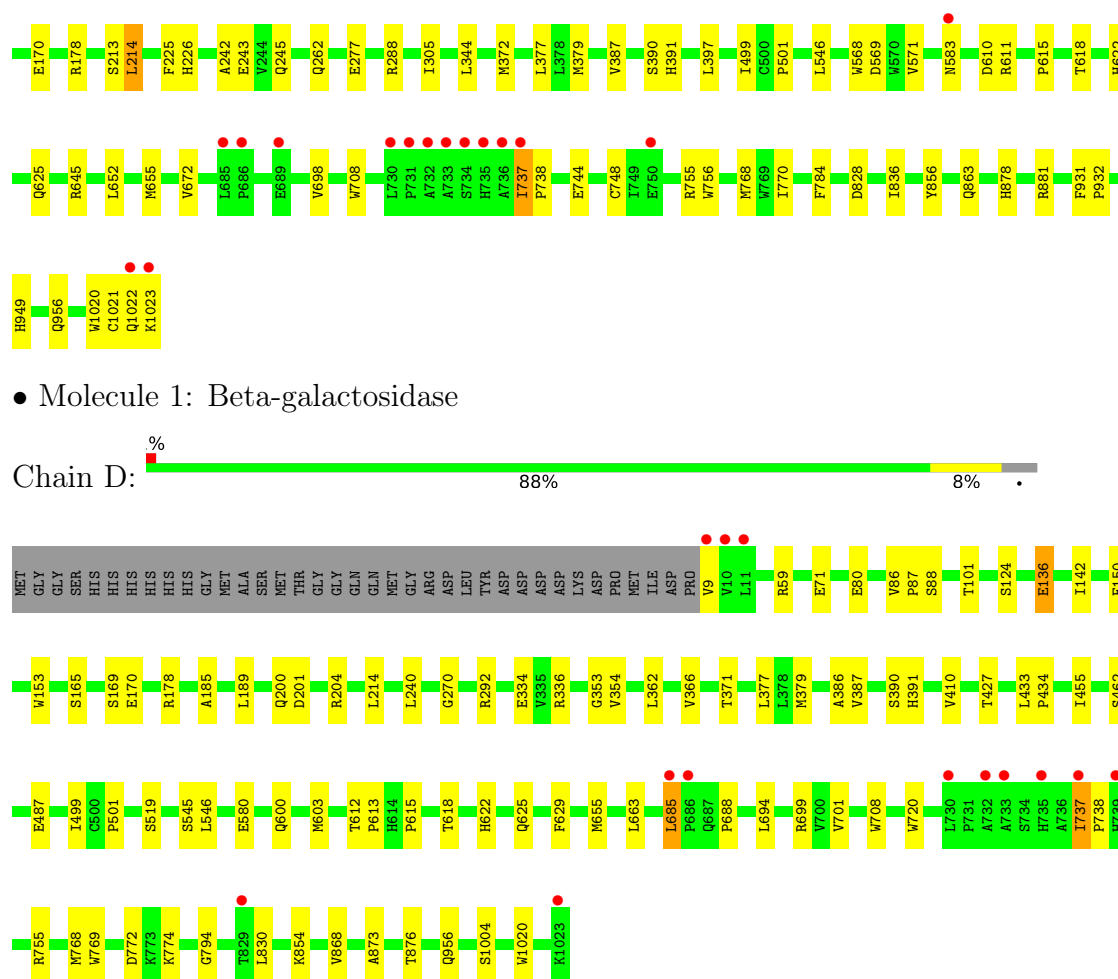


• Molecule 1: Beta-galactosidase



• 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, α , β , γ	152.60Å 164.69Å 203.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.05 – 2.50 46.05 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.05-2.50) 94.0 (46.05-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.5.0109, CNS	Depositor
R, R_{free}	0.174 , 0.237 0.178 , 0.237	Depositor DCC
R_{free} test set	2358 reflections (1.42%)	wwPDB-VP
Wilson B-factor (Å ²)	34.1	Xtrriage
Anisotropy	0.113	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	34381	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.13 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.4968e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: DMS, MG, 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	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.58	0/8398	0.78	3/11457 (0.0%)
1	B	0.59	0/8398	0.79	3/11457 (0.0%)
1	C	0.57	0/8398	0.77	1/11457 (0.0%)
1	D	0.57	0/8398	0.78	2/11457 (0.0%)
All	All	0.58	0/33592	0.78	9/45828 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	545	SER	N-CA-C	6.12	116.16	108.45
1	B	565	GLY	N-CA-C	5.86	119.16	110.42
1	A	690	SER	N-CA-C	5.83	115.80	108.45
1	B	233	ASP	N-CA-C	5.67	118.19	111.33
1	D	150	PHE	N-CA-C	5.35	117.22	108.76
1	B	150	PHE	N-CA-C	5.21	117.00	108.76
1	A	150	PHE	N-CA-C	5.17	116.92	108.76
1	A	833	ALA	N-CA-C	5.08	115.24	108.38
1	C	583	ASN	N-CA-C	5.01	116.88	109.62

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	8156	0	7754	28	0
1	B	8156	0	7754	30	0
1	C	8156	0	7754	36	0
1	D	8156	0	7754	39	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	3	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	3	0	0	0	0
4	A	28	0	42	1	0
4	B	32	0	48	1	0
4	C	28	0	42	1	0
4	D	40	0	60	3	0
5	A	427	0	0	1	0
5	B	403	0	0	3	0
5	C	389	0	0	2	0
5	D	392	0	0	5	0
All	All	34381	0	31208	131	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 (131) 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:D:142:ILE:HG12	1:D:170:GLU:HG2	1.60	0.84
1:A:142:ILE:HG12	1:A:170:GLU:HG2	1.67	0.76
1:C:652:LEU:HD11	1:C:698:VAL:HB	1.69	0.74
1:B:142:ILE:HG12	1:B:170:GLU:HG2	1.69	0.74
1:C:142:ILE:HG12	1:C:170:GLU:HG2	1.72	0.70
1:C:379:MET:HE1	1:C:387:VAL:HB	1.73	0.70
1:D:292:ARG:O	4:D:8010:DMS:H12	1.93	0.69
1:B:91:GLN:HG3	1:B:96:ASP:OD1	1.96	0.65
1:C:54:LEU:HD11	1:C:214:LEU:HD13	1.81	0.63
1:A:379:MET:HE1	1:A:387:VAL:HB	1.81	0.63
1:B:755:ARG:HB3	1:B:769:TRP:HB2	1.80	0.62
1:A:245:GLN:HG2	1:A:288:ARG:HG2	1.82	0.61
1:C:770:ILE:HD11	1:C:1022:GLN:HG2	1.83	0.60
1:C:305:ILE:HD11	1:C:645:ARG:HB3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:245:GLN:HG2	1:C:288:ARG:HG2	1.83	0.58
1:D:379:MET:HE1	1:D:387:VAL:HB	1.87	0.57
1:B:354:VAL:CG1	1:B:379:MET:HE3	2.34	0.57
1:A:9:VAL:HG12	1:D:9:VAL:HG12	1.87	0.57
1:C:92:MET:HE2	1:C:92:MET:HA	1.87	0.56
1:B:245:GLN:HG2	1:B:288:ARG:HG2	1.87	0.55
1:D:377:LEU:HD22	1:D:708:TRP:HA	1.89	0.55
1:D:487:GLU:HB3	5:D:4253:HOH:O	2.07	0.55
1:C:863:GLN:HG2	1:C:1021:CYS:HB3	1.90	0.53
1:B:360:HIS:CE1	1:B:362:LEU:HB2	2.43	0.53
1:A:788:PRO:HD2	1:A:968:MET:HB2	1.91	0.53
1:C:499:ILE:HG22	1:C:501:PRO:HD3	1.91	0.53
1:D:499:ILE:HG22	1:D:501:PRO:HD3	1.90	0.52
1:A:873:ALA:O	1:A:876:THR:HG22	2.09	0.52
1:D:88:SER:HA	1:D:366:VAL:HG21	1.92	0.51
1:A:887:GLN:NE2	1:A:980:GLU:O	2.44	0.51
1:B:379:MET:HE1	1:B:387:VAL:HB	1.91	0.51
1:D:270:GLY:HA3	4:D:8010:DMS:H13	1.91	0.50
4:B:8003:DMS:H13	5:B:4087:HOH:O	2.10	0.50
1:D:427:THR:HG21	1:D:462:SER:HB3	1.93	0.50
1:C:91:GLN:HG3	1:C:96:ASP:OD1	2.12	0.50
1:B:487:GLU:HB3	5:B:4039:HOH:O	2.11	0.50
1:D:629:PHE:HB3	1:D:720:TRP:HH2	1.77	0.49
1:D:688:PRO:HG3	1:D:694:LEU:HD11	1.95	0.49
1:C:84:VAL:HG12	4:C:8006:DMS:H23	1.95	0.49
1:B:127:PHE:CE1	1:B:184:LEU:HG	2.48	0.48
1:A:147:ASN:HA	1:A:148:SER:HA	1.62	0.48
1:D:755:ARG:HB3	1:D:769:TRP:HB2	1.96	0.48
1:B:651:LEU:HD12	1:B:651:LEU:C	2.39	0.47
1:B:794:GLY:HA2	1:B:998:SER:O	2.15	0.47
1:A:830:LEU:HD22	1:B:828:ASP:HB3	1.97	0.47
1:B:521:LYS:HG2	1:B:559:TYR:CZ	2.49	0.47
1:C:91:GLN:HB3	1:C:98:PRO:HD3	1.95	0.47
1:D:410:VAL:HG22	1:D:455:ILE:HB	1.96	0.47
1:D:655:MET:HB2	5:D:4353:HOH:O	2.13	0.47
1:C:568:TRP:CD2	1:C:569:ASP:HB3	2.50	0.47
1:C:949:HIS:CD2	1:C:1020:TRP:HE1	2.33	0.46
1:D:615:PRO:O	1:D:618:THR:HG22	2.14	0.46
1:C:225:PHE:HA	1:C:243:GLU:O	2.15	0.46
1:D:240:LEU:C	1:D:240:LEU:HD23	2.41	0.46
1:A:427:THR:HG21	1:A:462:SER:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:TYR:CZ	1:A:602:CYS:HB3	2.51	0.46
1:A:200:GLN:HA	1:A:416:GLU:OE1	2.16	0.46
1:D:768:MET:HE1	1:D:1020:TRP:CZ2	2.51	0.45
1:D:136:GLU:HG2	5:D:4181:HOH:O	2.15	0.45
1:D:101:THR:HG23	1:D:204:ARG:NH2	2.31	0.45
1:D:622:HIS:NE2	4:D:8002:DMS:H13	2.32	0.44
1:D:353:GLY:HA2	1:D:386:ALA:O	2.18	0.44
1:A:114:VAL:HB	1:A:115:PRO:CD	2.47	0.44
1:C:784:PHE:HA	1:C:881:ARG:O	2.18	0.44
1:D:153:TRP:HB2	1:D:185:ALA:HB3	1.98	0.44
1:D:873:ALA:O	1:D:876:THR:HG22	2.18	0.44
1:A:763:GLY:HA3	1:A:822:LEU:HD13	1.99	0.44
1:C:931:PHE:HA	1:C:932:PRO:HD3	1.78	0.44
1:B:127:PHE:HE1	1:B:184:LEU:HG	1.81	0.44
1:B:615:PRO:O	1:B:618:THR:HG22	2.18	0.43
1:C:377:LEU:HD22	1:C:708:TRP:HA	2.00	0.43
1:D:200:GLN:HG2	1:D:391:HIS:HB2	1.99	0.43
1:D:737:ILE:HA	1:D:738:PRO:HD3	1.87	0.43
1:D:334:GLU:OE1	1:D:336:ARG:NH1	2.48	0.43
1:A:581:ASN:HB2	5:A:4192:HOH:O	2.18	0.43
1:B:782:ASP:HB2	1:B:842:TRP:CH2	2.53	0.43
1:C:828:ASP:HB3	1:D:830:LEU:HD22	2.00	0.43
1:B:745:MET:HE2	1:B:745:MET:HA	2.00	0.43
1:B:784:PHE:HA	1:B:881:ARG:O	2.19	0.43
1:C:836:ILE:HB	1:C:856:TYR:HB2	2.00	0.43
1:C:622:HIS:O	1:C:625:GLN:HG3	2.19	0.43
1:A:772:ASP:OD1	1:A:772:ASP:N	2.51	0.43
1:B:651:LEU:HD11	1:B:653:HIS:CE1	2.54	0.43
1:D:685:LEU:H	1:D:685:LEU:HG	1.42	0.43
1:D:854:LYS:HG2	1:D:868:VAL:HG22	2.01	0.43
1:A:390:SER:HA	1:A:391:HIS:HA	1.86	0.42
1:C:568:TRP:HA	1:C:569:ASP:HA	1.88	0.42
1:C:768:MET:HE1	1:C:1020:TRP:CZ2	2.54	0.42
1:D:433:LEU:HB3	1:D:434:PRO:HD3	2.01	0.42
1:D:86:VAL:HG13	1:D:87:PRO:HA	2.01	0.42
1:A:129:VAL:HG21	1:A:177:LEU:HD22	2.02	0.42
1:C:737:ILE:HA	1:C:738:PRO:HD3	1.92	0.42
1:C:878:HIS:HD2	5:C:4072:HOH:O	2.02	0.42
1:A:864:MET:HE3	1:A:866:ILE:HD11	2.01	0.42
1:C:147:ASN:HA	1:C:148:SER:HA	1.72	0.42
1:B:713:HIS:HE1	5:B:4262:HOH:O	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:600:GLN:HB2	1:D:603:MET:HE2	2.01	0.42
1:B:883:GLY:HA3	1:B:987:ASP:HA	2.01	0.42
1:D:59:ARG:HB2	1:D:124:SER:OG	2.20	0.42
1:A:36:TRP:HH2	4:A:8004:DMS:H23	1.85	0.42
1:A:200:GLN:HG3	1:A:416:GLU:OE1	2.20	0.42
1:B:100:TYR:CE1	1:B:602:CYS:HB3	2.54	0.42
1:A:377:LEU:HD22	1:A:708:TRP:HA	2.02	0.41
1:C:615:PRO:O	1:C:618:THR:HG22	2.20	0.41
1:A:505:ARG:O	1:A:519:SER:HA	2.21	0.41
1:B:88:SER:HA	1:B:366:VAL:HG21	2.02	0.41
1:C:379:MET:CE	1:C:387:VAL:HB	2.47	0.41
1:C:610:ASP:O	1:C:611:ARG:HB2	2.20	0.41
1:C:748:CYS:HA	1:C:756:TRP:O	2.21	0.41
1:D:625:GLN:NE2	5:D:4062:HOH:O	2.52	0.41
1:A:240:LEU:C	1:A:240:LEU:HD23	2.46	0.41
1:B:730:LEU:H	1:B:730:LEU:HD12	1.86	0.41
1:C:226:HIS:O	1:C:242:ALA:HA	2.20	0.41
1:C:277:GLU:HG2	5:C:4075:HOH:O	2.21	0.41
1:A:147:ASN:O	1:A:206:SER:HB2	2.21	0.41
1:A:610:ASP:O	1:A:611:ARG:HB2	2.21	0.41
1:A:499:ILE:HG22	1:A:501:PRO:HD3	2.02	0.41
1:B:568:TRP:CD2	1:B:569:ASP:HB3	2.56	0.41
1:C:372:MET:CE	1:C:397:LEU:HD23	2.51	0.41
1:C:390:SER:HA	1:C:391:HIS:HA	1.82	0.41
1:D:354:VAL:CG1	1:D:379:MET:HE3	2.51	0.41
1:B:369:GLU:HA	1:B:372:MET:HE3	2.03	0.41
1:A:883:GLY:HA3	1:A:987:ASP:HA	2.03	0.40
1:B:100:TYR:CZ	1:B:602:CYS:HB3	2.56	0.40
1:D:390:SER:HA	1:D:391:HIS:HA	1.75	0.40
1:B:60:PHE:HA	1:B:122:CYS:O	2.21	0.40
1:C:756:TRP:CD1	1:C:768:MET:HG2	2.56	0.40
1:B:126:THR:HA	1:B:182:ASN:O	2.21	0.40
1:D:794:GLY:HA3	5:D:4071:HOH:O	2.20	0.40
1:B:542:MET:HE3	1:B:601:PHE:HA	2.03	0.40
1:D:612:THR:HA	1:D:613:PRO:HD3	1.97	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	1013/1052 (96%)	974 (96%)	39 (4%)	0	100	100
1	B	1013/1052 (96%)	967 (96%)	46 (4%)	0	100	100
1	C	1013/1052 (96%)	975 (96%)	38 (4%)	0	100	100
1	D	1013/1052 (96%)	976 (96%)	36 (4%)	1 (0%)	48	68
All	All	4052/4208 (96%)	3892 (96%)	159 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	201	ASP

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	868/898 (97%)	852 (98%)	16 (2%)	51	77
1	B	868/898 (97%)	853 (98%)	15 (2%)	53	78
1	C	868/898 (97%)	852 (98%)	16 (2%)	51	77
1	D	868/898 (97%)	846 (98%)	22 (2%)	42	69
All	All	3472/3592 (97%)	3403 (98%)	69 (2%)	48	75

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	251	ARG
1	A	319	ASP
1	A	344	LEU
1	A	519	SER
1	A	546	LEU
1	A	599	ARG
1	A	651	LEU
1	A	652	LEU
1	A	672	VAL
1	A	684	GLU
1	A	744	GLU
1	A	750	GLU
1	A	772	ASP
1	A	778	THR
1	A	910	LEU
1	A	956	GLN
1	B	101	THR
1	B	265	THR
1	B	291	LEU
1	B	362	LEU
1	B	519	SER
1	B	546	LEU
1	B	651	LEU
1	B	652	LEU
1	B	672	VAL
1	B	684	GLU
1	B	689	GLU
1	B	737	ILE
1	B	772	ASP
1	B	910	LEU
1	B	956	GLN
1	C	116	THR
1	C	130	ASP
1	C	178	ARG
1	C	213	SER
1	C	214	LEU
1	C	262	GLN
1	C	344	LEU
1	C	546	LEU
1	C	571	VAL
1	C	655	MET
1	C	672	VAL
1	C	737	ILE

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Mol	Chain	Res	Type
1	C	744	GLU
1	C	755	ARG
1	C	956	GLN
1	C	1023	LYS
1	D	71	GLU
1	D	80	GLU
1	D	136	GLU
1	D	165	SER
1	D	169	SER
1	D	178	ARG
1	D	189	LEU
1	D	214	LEU
1	D	362	LEU
1	D	371	THR
1	D	519	SER
1	D	546	LEU
1	D	580	GLU
1	D	663	LEU
1	D	685	LEU
1	D	699	ARG
1	D	701	VAL
1	D	737	ILE
1	D	772	ASP
1	D	774	LYS
1	D	956	GLN
1	D	1004	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	307	ASN
1	A	634	GLN
1	A	653	HIS
1	A	757	GLN
1	A	761	GLN
1	A	844	HIS
1	A	945	ASN
1	A	950	GLN
1	A	974	HIS
1	B	138	GLN
1	B	221	GLN

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Mol	Chain	Res	Type
1	B	394	ASN
1	B	485	GLN
1	B	623	GLN
1	B	624	GLN
1	B	653	HIS
1	B	739	HIS
1	B	757	GLN
1	B	878	HIS
1	B	887	GLN
1	C	12	GLN
1	C	262	GLN
1	C	307	ASN
1	C	445	GLN
1	C	783	GLN
1	C	896	ASN
1	C	945	ASN
1	C	974	HIS
1	C	985	ASN
1	C	1022	GLN
1	D	163	GLN
1	D	624	GLN
1	D	634	GLN
1	D	653	HIS
1	D	687	GLN
1	D	757	GLN
1	D	945	ASN
1	D	950	GLN
1	D	977	HIS
1	D	985	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 50 ligands modelled in this entry, 18 are monoatomic - leaving 32 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
4	DMS	B	8007	-	3,3,3	2.69	1 (33%)	3,3,3	0.55	0
4	DMS	A	8006	-	3,3,3	2.64	1 (33%)	3,3,3	0.36	0
4	DMS	C	8001	-	3,3,3	2.60	1 (33%)	3,3,3	0.60	0
4	DMS	C	8003	-	3,3,3	2.75	1 (33%)	3,3,3	0.52	0
4	DMS	C	8005	-	3,3,3	2.67	1 (33%)	3,3,3	0.47	0
4	DMS	B	8008	-	3,3,3	2.73	1 (33%)	3,3,3	0.50	0
4	DMS	D	8002	-	3,3,3	2.66	1 (33%)	3,3,3	0.69	0
4	DMS	D	8006	-	3,3,3	2.73	1 (33%)	3,3,3	0.66	0
4	DMS	D	8010	-	3,3,3	2.68	1 (33%)	3,3,3	0.41	0
4	DMS	B	8004	-	3,3,3	2.74	1 (33%)	3,3,3	0.74	0
4	DMS	C	8004	-	3,3,3	2.77	1 (33%)	3,3,3	0.78	0
4	DMS	D	8005	-	3,3,3	2.72	1 (33%)	3,3,3	0.63	0
4	DMS	B	8003	-	3,3,3	2.70	1 (33%)	3,3,3	0.58	0
4	DMS	A	8004	-	3,3,3	2.75	1 (33%)	3,3,3	0.66	0
4	DMS	D	8008	-	3,3,3	2.74	1 (33%)	3,3,3	0.60	0
4	DMS	A	8003	-	3,3,3	2.72	1 (33%)	3,3,3	0.65	0
4	DMS	A	8001	-	3,3,3	2.64	1 (33%)	3,3,3	0.81	0
4	DMS	A	8005	-	3,3,3	2.66	1 (33%)	3,3,3	0.62	0
4	DMS	A	8002	-	3,3,3	2.73	1 (33%)	3,3,3	0.47	0
4	DMS	D	8004	-	3,3,3	2.77	1 (33%)	3,3,3	0.71	0
4	DMS	B	8002	-	3,3,3	2.65	1 (33%)	3,3,3	0.76	0
4	DMS	D	8007	-	3,3,3	2.68	1 (33%)	3,3,3	0.68	0
4	DMS	C	8002	-	3,3,3	2.69	1 (33%)	3,3,3	0.60	0
4	DMS	D	8009	-	3,3,3	2.73	1 (33%)	3,3,3	0.55	0
4	DMS	D	8001	-	3,3,3	2.68	1 (33%)	3,3,3	0.57	0
4	DMS	B	8001	-	3,3,3	2.62	1 (33%)	3,3,3	0.74	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DMS	D	8003	-	3,3,3	2.73	1 (33%)	3,3,3	0.37	0
4	DMS	C	8007	-	3,3,3	2.69	1 (33%)	3,3,3	0.47	0
4	DMS	B	8006	-	3,3,3	2.77	1 (33%)	3,3,3	0.55	0
4	DMS	B	8005	-	3,3,3	2.66	1 (33%)	3,3,3	0.49	0
4	DMS	A	8007	-	3,3,3	2.69	1 (33%)	3,3,3	0.55	0
4	DMS	C	8006	-	3,3,3	2.72	1 (33%)	3,3,3	0.54	0

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	8004	DMS	O-S	4.64	1.80	1.50
4	B	8006	DMS	O-S	4.63	1.80	1.50
4	C	8003	DMS	O-S	4.62	1.80	1.50
4	A	8004	DMS	O-S	4.62	1.80	1.50
4	C	8004	DMS	O-S	4.61	1.80	1.50
4	A	8002	DMS	O-S	4.61	1.80	1.50
4	B	8004	DMS	O-S	4.60	1.80	1.50
4	D	8008	DMS	O-S	4.59	1.80	1.50
4	B	8008	DMS	O-S	4.59	1.80	1.50
4	D	8003	DMS	O-S	4.59	1.80	1.50
4	D	8006	DMS	O-S	4.57	1.80	1.50
4	D	8005	DMS	O-S	4.57	1.80	1.50
4	D	8009	DMS	O-S	4.56	1.80	1.50
4	C	8006	DMS	O-S	4.56	1.80	1.50
4	A	8003	DMS	O-S	4.55	1.80	1.50
4	B	8003	DMS	O-S	4.53	1.80	1.50
4	D	8010	DMS	O-S	4.52	1.80	1.50
4	C	8002	DMS	O-S	4.51	1.80	1.50
4	C	8007	DMS	O-S	4.51	1.80	1.50
4	A	8007	DMS	O-S	4.50	1.79	1.50
4	B	8007	DMS	O-S	4.49	1.79	1.50
4	D	8001	DMS	O-S	4.49	1.79	1.50
4	D	8007	DMS	O-S	4.48	1.79	1.50
4	C	8005	DMS	O-S	4.47	1.79	1.50
4	D	8002	DMS	O-S	4.46	1.79	1.50
4	B	8002	DMS	O-S	4.44	1.79	1.50
4	B	8005	DMS	O-S	4.44	1.79	1.50
4	A	8005	DMS	O-S	4.44	1.79	1.50
4	A	8001	DMS	O-S	4.43	1.79	1.50
4	A	8006	DMS	O-S	4.42	1.79	1.50
4	B	8001	DMS	O-S	4.40	1.79	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	8001	DMS	O-S	4.34	1.78	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	8002	DMS	1	0
4	D	8010	DMS	2	0
4	B	8003	DMS	1	0
4	A	8004	DMS	1	0
4	C	8006	DMS	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	1015/1052 (96%)	-0.25	13 (1%) 75 71	28, 41, 64, 96	0
1	B	1015/1052 (96%)	-0.18	23 (2%) 61 57	25, 42, 66, 109	0
1	C	1015/1052 (96%)	-0.18	17 (1%) 69 65	27, 42, 66, 99	0
1	D	1015/1052 (96%)	-0.20	13 (1%) 75 71	26, 42, 66, 101	0
All	All	4060/4208 (96%)	-0.20	66 (1%) 70 67	25, 42, 66, 109	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	LEU	3.9
1	C	733	ALA	3.8
1	D	685	LEU	3.8
1	B	730	LEU	3.7
1	C	10	VAL	3.5
1	B	761	GLN	3.5
1	A	10	VAL	3.4
1	B	11	LEU	3.4
1	A	732	ALA	3.4
1	B	733	ALA	3.4
1	B	731	PRO	3.3
1	C	730	LEU	3.3
1	D	11	LEU	3.3
1	B	10	VAL	3.3
1	A	686	PRO	3.2
1	B	736	ALA	3.2
1	A	730	LEU	3.1
1	C	11	LEU	3.0
1	A	9	VAL	3.0
1	B	734	SER	3.0
1	B	739	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	731	PRO	2.9
1	B	732	ALA	2.8
1	C	1022	GLN	2.8
1	D	686	PRO	2.8
1	A	772	ASP	2.8
1	C	737	ILE	2.8
1	C	689	GLU	2.8
1	B	735	HIS	2.8
1	A	733	ALA	2.7
1	B	755	ARG	2.7
1	C	734	SER	2.7
1	B	770	ILE	2.7
1	B	9	VAL	2.6
1	D	735	HIS	2.6
1	C	750	GLU	2.5
1	C	1023	LYS	2.4
1	D	733	ALA	2.4
1	D	730	LEU	2.4
1	D	739	HIS	2.4
1	C	686	PRO	2.4
1	C	735	HIS	2.4
1	B	774	LYS	2.4
1	D	732	ALA	2.4
1	C	685	LEU	2.3
1	C	736	ALA	2.3
1	D	10	VAL	2.3
1	D	737	ILE	2.3
1	B	831	ALA	2.2
1	C	583	ASN	2.2
1	A	735	HIS	2.2
1	D	9	VAL	2.1
1	B	729	THR	2.1
1	B	686	PRO	2.1
1	D	1023	LYS	2.1
1	A	655	MET	2.1
1	C	732	ALA	2.1
1	B	690	SER	2.1
1	A	731	PRO	2.1
1	D	829	THR	2.1
1	A	685	LEU	2.0
1	B	684	GLU	2.0
1	B	689	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	772	ASP	2.0
1	B	688	PRO	2.0
1	A	734	SER	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 oligosaccharides 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
4	DMS	C	8007	4/4	0.85	0.20	71,73,74,74	0
4	DMS	D	8004	4/4	0.86	0.20	76,77,77,80	0
4	DMS	C	8004	4/4	0.87	0.20	75,76,76,78	0
4	DMS	A	8007	4/4	0.87	0.21	74,75,78,78	0
4	DMS	B	8004	4/4	0.87	0.20	67,70,71,71	0
4	DMS	A	8004	4/4	0.88	0.23	77,78,79,79	0
4	DMS	A	8006	4/4	0.89	0.21	65,67,69,69	0
4	DMS	C	8006	4/4	0.90	0.20	82,82,82,83	0
4	DMS	D	8006	4/4	0.90	0.19	78,79,80,81	0
4	DMS	B	8005	4/4	0.91	0.20	69,71,72,72	0
4	DMS	B	8006	4/4	0.92	0.14	62,64,64,67	0
4	DMS	D	8009	4/4	0.92	0.17	72,73,74,76	0
4	DMS	B	8007	4/4	0.93	0.16	72,72,73,74	0
4	DMS	C	8005	4/4	0.93	0.14	67,68,68,69	0
4	DMS	D	8008	4/4	0.93	0.15	61,61,62,64	0
4	DMS	D	8002	4/4	0.93	0.19	68,68,69,71	0
3	NA	D	3102	1/1	0.94	0.06	35,35,35,35	0
3	NA	D	3103	1/1	0.94	0.09	51,51,51,51	0
4	DMS	A	8003	4/4	0.94	0.14	52,53,54,55	0
3	NA	A	3102	1/1	0.94	0.06	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	DMS	A	8005	4/4	0.94	0.16	55,58,59,60	0
4	DMS	D	8005	4/4	0.94	0.13	75,75,76,78	0
4	DMS	C	8001	4/4	0.94	0.14	50,52,54,55	0
4	DMS	C	8003	4/4	0.94	0.13	64,65,65,66	0
3	NA	C	3101	1/1	0.94	0.10	44,44,44,44	0
4	DMS	A	8001	4/4	0.95	0.16	52,54,55,56	0
4	DMS	B	8008	4/4	0.95	0.11	48,49,50,51	0
2	MG	B	3002	1/1	0.95	0.12	46,46,46,46	0
4	DMS	C	8002	4/4	0.95	0.13	56,56,58,59	0
4	DMS	B	8002	4/4	0.95	0.13	51,52,53,53	0
4	DMS	D	8010	4/4	0.95	0.15	54,55,56,61	0
3	NA	A	3101	1/1	0.96	0.05	41,41,41,41	0
4	DMS	B	8003	4/4	0.96	0.12	56,57,57,58	0
3	NA	C	3102	1/1	0.96	0.04	33,33,33,33	0
4	DMS	A	8002	4/4	0.96	0.13	48,49,50,52	0
2	MG	A	3002	1/1	0.96	0.05	39,39,39,39	0
4	DMS	D	8003	4/4	0.96	0.12	55,55,57,57	0
2	MG	C	3002	1/1	0.97	0.10	44,44,44,44	0
4	DMS	D	8001	4/4	0.97	0.09	51,51,51,53	0
4	DMS	D	8007	4/4	0.97	0.11	62,65,65,67	0
2	MG	D	3002	1/1	0.97	0.06	45,45,45,45	0
4	DMS	B	8001	4/4	0.97	0.11	47,48,50,51	0
3	NA	B	3102	1/1	0.97	0.04	31,31,31,31	0
3	NA	A	3103	1/1	0.98	0.07	41,41,41,41	0
3	NA	D	3101	1/1	0.98	0.07	43,43,43,43	0
3	NA	B	3101	1/1	0.98	0.07	44,44,44,44	0
2	MG	D	3001	1/1	0.99	0.04	39,39,39,39	0
2	MG	A	3001	1/1	0.99	0.03	39,39,39,39	0
2	MG	C	3001	1/1	0.99	0.03	40,40,40,40	0
2	MG	B	3001	1/1	0.99	0.03	39,39,39,39	0

6.5 Other polymers ⓘ

There are no such residues in this entry.