



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:41 PM UTC

PDB ID : 3VD3 / pdb_00003vd3
Title : E. coli (lacZ) beta-galactosidase (N460D)
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 : 2.80 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

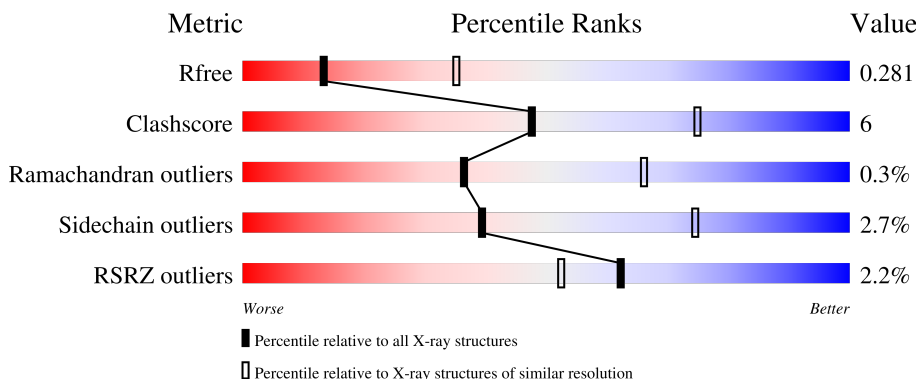
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	
1	B	1052	
1	C	1052	
1	D	1052	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 33130 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	1011	Total	C	N	O	S	0	0	0
			8126	5138	1439	1511	38			
1	B	1011	Total	C	N	O	S	0	0	0
			8126	5138	1439	1511	38			
1	C	1011	Total	C	N	O	S	0	0	0
			8126	5138	1439	1511	38			
1	D	1011	Total	C	N	O	S	0	0	0
			8126	5138	1439	1511	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	ASP	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	ASP	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	ASP	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	ASP	ASN	engineered mutation	UNP P00722

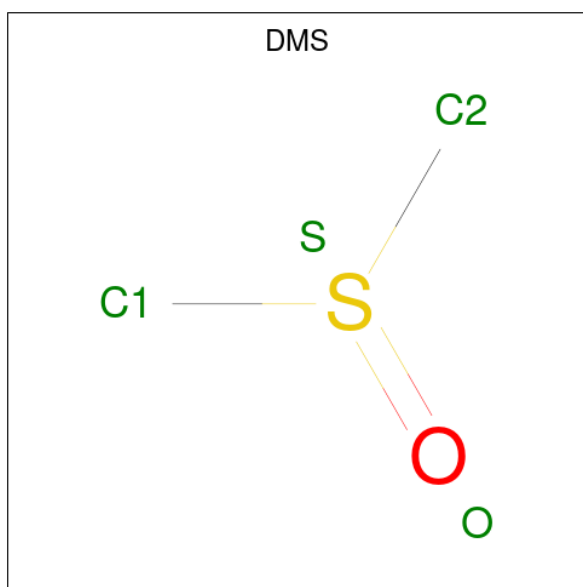
- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
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	4	Total	Na	0	0
			4	4		
3	C	3	Total	Na	0	0
			3	3		
3	D	2	Total	Na	0	0
			2	2		

- 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	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		

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

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

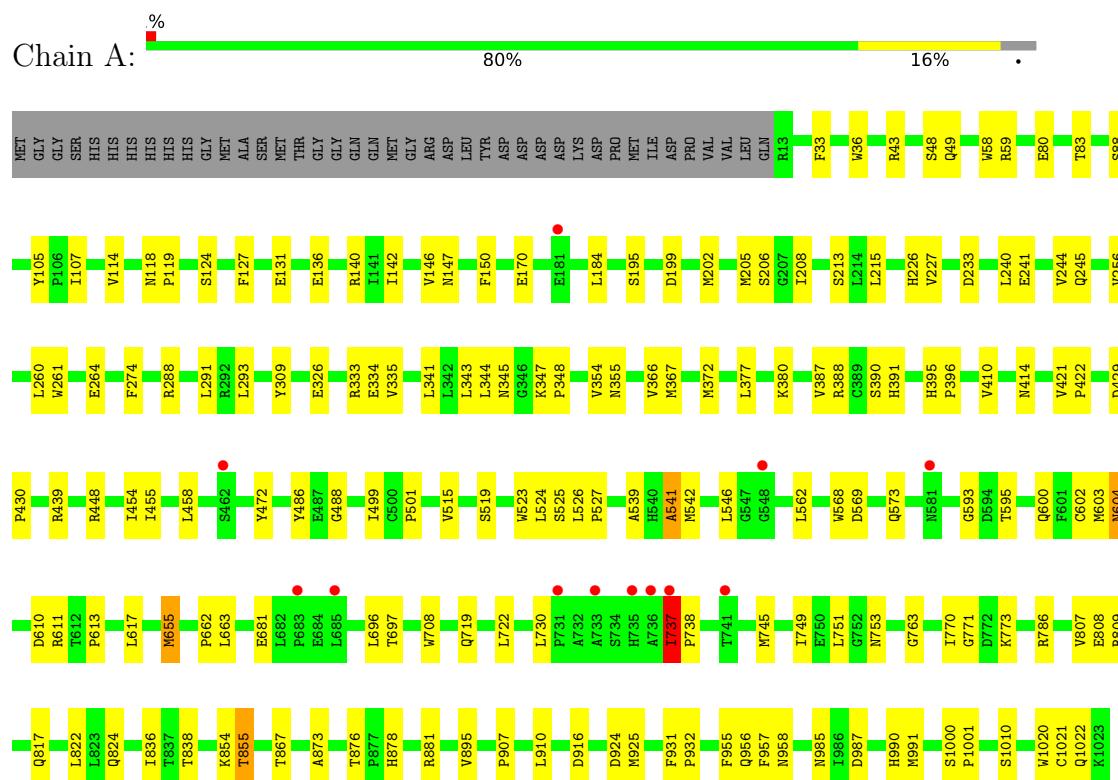
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	112	Total O 112 112	0	0
5	B	110	Total O 110 110	0	0
5	C	99	Total O 99 99	0	0
5	D	106	Total O 106 106	0	0

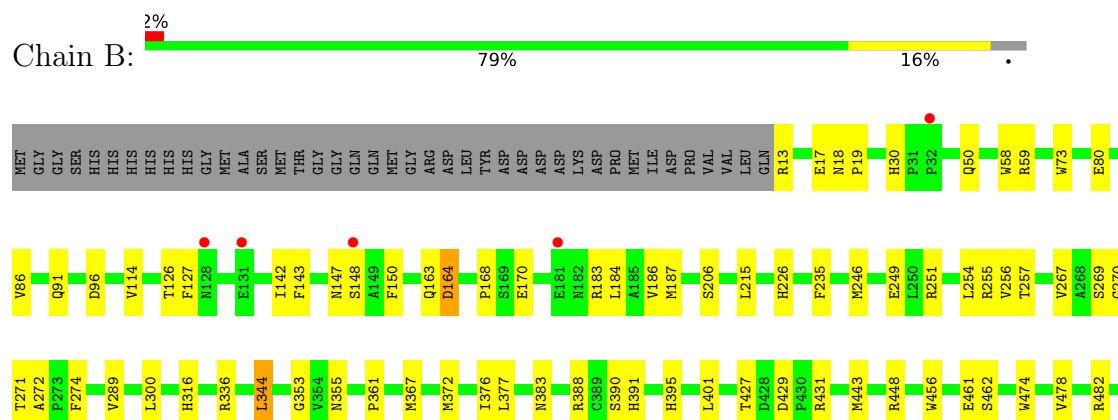
3 Residue-property plots [i](#)

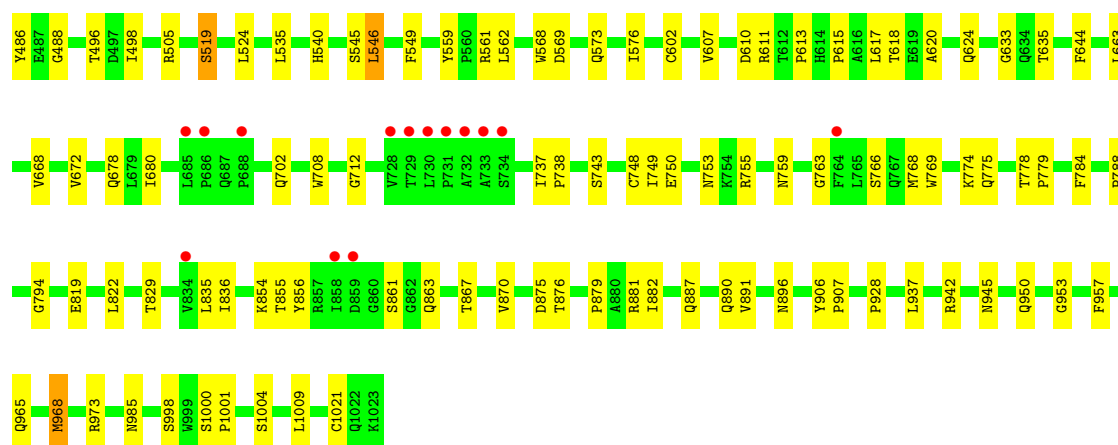
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

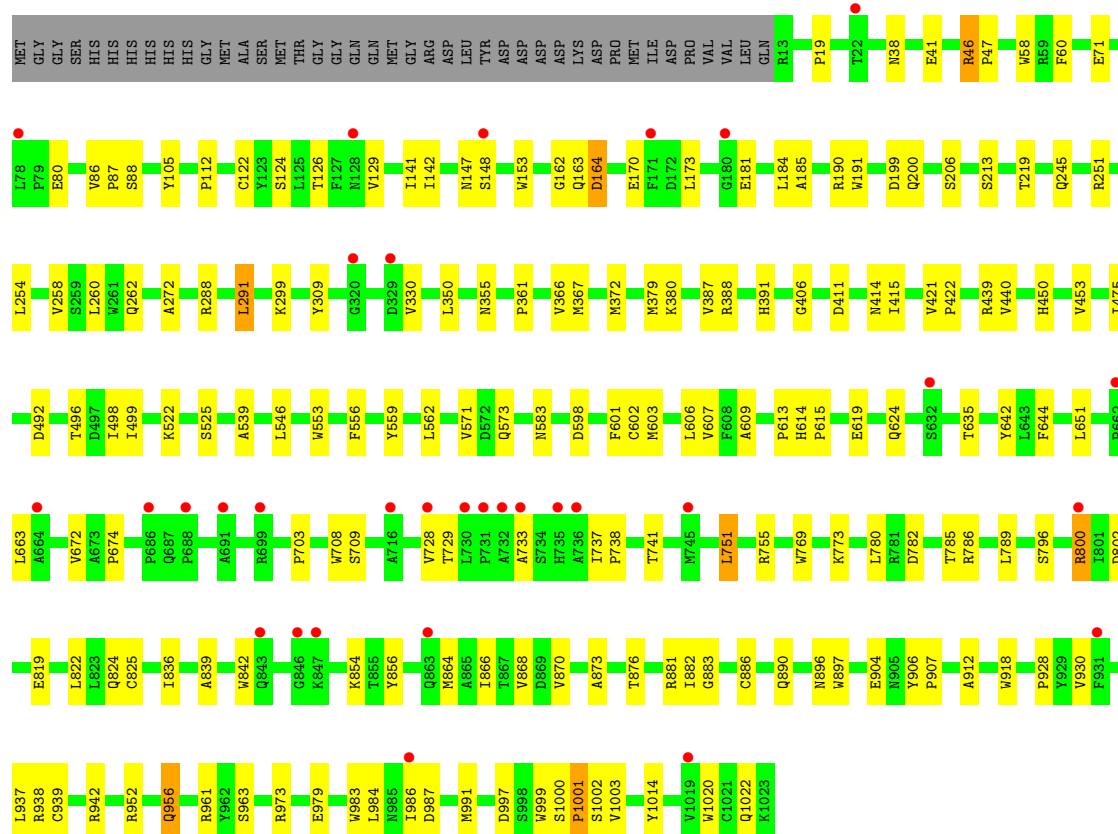
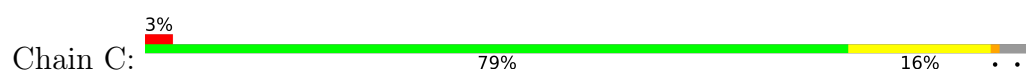


• Molecule 1: Beta-galactosidase

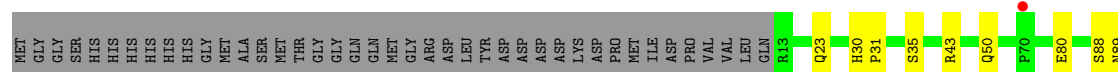
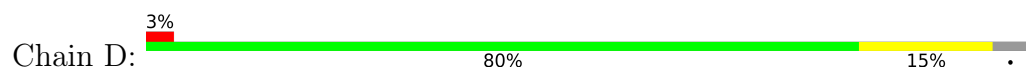


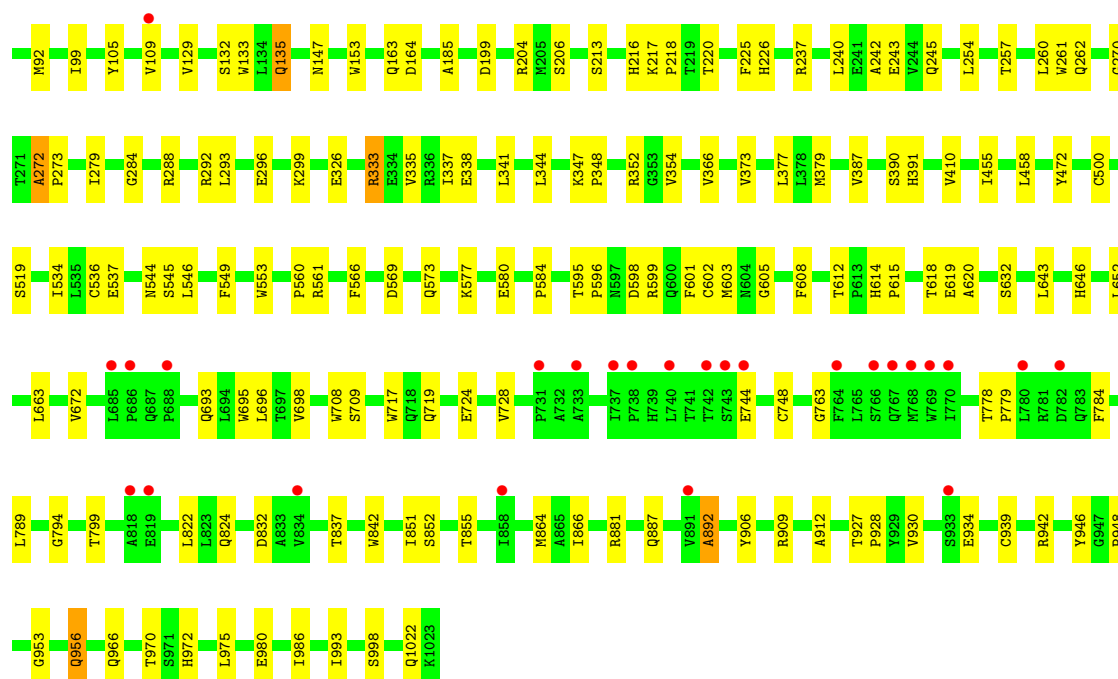


• Molecule 1: Beta-galactosidase



• Molecule 1: Beta-galactosidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	127.14Å 151.65Å 131.33Å 90.00° 103.47° 90.00°	Depositor
Resolution (Å)	57.25 – 2.80 57.25 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (57.25-2.80) 88.0 (57.25-2.80)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0109, CNS	Depositor
R, R_{free}	0.200 , 0.282 0.201 , 0.281	Depositor DCC
R_{free} test set	5235 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 19.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.028 for l,-k,h	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	33130	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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: 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	0.53	1/8368 (0.0%)	0.80	3/11415 (0.0%)
1	B	0.51	0/8368	0.80	2/11415 (0.0%)
1	C	0.52	0/8368	0.80	7/11415 (0.1%)
1	D	0.54	0/8368	0.79	3/11415 (0.0%)
All	All	0.53	1/33472 (0.0%)	0.80	15/45660 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	737	ILE	CA-CB	5.74	1.58	1.54

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	272	ALA	N-CA-C	7.00	114.52	108.22
1	D	272	ALA	N-CA-C	6.60	114.07	108.13
1	A	233	ASP	N-CA-C	5.79	119.60	112.54
1	B	353	GLY	N-CA-C	5.75	119.54	111.14
1	C	571	VAL	N-CA-C	5.60	116.35	108.17
1	C	162	GLY	N-CA-C	5.40	120.03	110.80
1	C	272	ALA	N-CA-C	5.30	112.99	108.22
1	A	955	PHE	N-CA-C	5.30	116.71	109.18
1	C	450	HIS	CA-C-N	5.22	124.88	119.56
1	C	450	HIS	C-N-CA	5.22	124.88	119.56
1	D	892	ALA	N-CA-C	5.14	117.09	109.69
1	D	599	ARG	N-CA-C	5.14	116.12	108.31
1	A	593	GLY	N-CA-C	-5.10	108.58	115.36
1	C	583	ASN	CA-C-N	5.06	125.85	120.13
1	C	583	ASN	C-N-CA	5.06	125.85	120.13

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	8126	0	7714	91	0
1	B	8126	0	7714	93	0
1	C	8126	0	7714	88	0
1	D	8126	0	7714	90	0
2	A	1	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	4	0	0	0	0
3	C	3	0	0	0	0
3	D	2	0	0	0	0
4	A	48	0	72	2	0
4	B	44	0	66	2	0
4	C	48	0	72	3	0
4	D	40	0	60	3	0
5	A	112	0	0	6	0
5	B	110	0	0	3	0
5	C	99	0	0	1	0
5	D	106	0	0	6	0
All	All	33130	0	31126	358	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 6.

All (358) 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:573:GLN:HB2	1:B:602:CYS:O	1.82	0.79
1:A:367:MET:HB3	1:A:372:MET:HE3	1.66	0.77
1:A:245:GLN:HG2	1:A:288:ARG:HG2	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:832:ASP:HB3	5:D:4075:HOH:O	1.86	0.75
1:C:836:ILE:HB	1:C:856:TYR:HB2	1.68	0.75
1:C:642:TYR:OH	4:C:8002:DMS:H22	1.88	0.73
1:A:202:MET:HE2	5:A:4004:HOH:O	1.89	0.72
1:A:809:ARG:HD2	5:A:4045:HOH:O	1.91	0.71
1:B:127:PHE:HE2	1:B:184:LEU:HG	1.56	0.70
1:A:105:TYR:CE2	1:A:199:ASP:HB2	2.26	0.70
1:A:88:SER:HA	1:A:366:VAL:HG21	1.74	0.70
1:D:789:LEU:HD11	1:D:993:ILE:HG22	1.74	0.69
1:A:244:VAL:HG21	1:A:256:VAL:HG11	1.75	0.68
1:D:88:SER:HA	1:D:366:VAL:HG21	1.76	0.68
1:A:786:ARG:NH2	1:A:991:MET:HE2	2.09	0.67
1:D:352:ARG:HG2	1:D:553:TRP:CH2	2.30	0.67
1:B:945:ASN:OD1	1:B:950:GLN:HG3	1.95	0.66
1:A:573:GLN:HB2	1:A:602:CYS:O	1.96	0.66
1:B:127:PHE:CE2	1:B:184:LEU:HG	2.31	0.65
1:A:873:ALA:O	1:A:876:THR:HG22	1.96	0.64
1:B:854:LYS:HA	1:B:867:THR:O	1.98	0.64
1:A:136:GLU:HB2	5:A:4033:HOH:O	1.97	0.64
1:D:619:GLU:HA	1:D:912:ALA:HB2	1.80	0.63
1:D:892:ALA:HB3	1:D:946:TYR:CE1	2.35	0.62
1:D:333:ARG:NH1	4:D:8001:DMS:H12	2.15	0.62
1:A:410:VAL:HG22	1:A:455:ILE:HB	1.81	0.61
1:A:895:VAL:HG21	1:A:925:MET:HG3	1.82	0.61
1:D:763:GLY:HA3	1:D:822:LEU:HD13	1.80	0.61
1:C:780:LEU:HD13	1:C:1020:TRP:HZ3	1.65	0.61
1:A:749:ILE:HD11	1:A:836:ILE:HD11	1.81	0.61
1:B:784:PHE:HA	1:B:881:ARG:O	2.00	0.61
1:D:939:CYS:HA	1:D:956:GLN:HB3	1.82	0.61
1:A:377:LEU:HD22	1:A:708:TRP:HA	1.83	0.60
1:D:948:PRO:HB2	1:D:1022:GLN:HG2	1.82	0.60
1:C:651:LEU:HD23	1:C:703:PRO:HG3	1.82	0.60
1:C:782:ASP:HB2	1:C:842:TRP:CH2	2.36	0.59
1:A:36:TRP:CD1	1:A:48:SER:HB2	2.37	0.59
1:C:200:GLN:HG2	1:C:391:HIS:HB2	1.84	0.59
1:C:559:TYR:HB2	1:C:562:LEU:HD12	1.85	0.59
1:B:367:MET:HB3	1:B:372:MET:HE2	1.85	0.58
1:A:458:LEU:HD11	1:A:472:TYR:HB2	1.86	0.58
1:B:73:TRP:O	1:B:183:ARG:NH1	2.27	0.58
1:A:525:SER:O	1:B:561:ARG:HD3	2.03	0.58
1:B:59:ARG:HG2	4:B:8011:DMS:H13	1.84	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:573:GLN:HB2	1:C:602:CYS:O	2.05	0.56
1:B:163:GLN:O	1:B:164:ASP:CB	2.53	0.56
1:A:499:ILE:HG22	1:A:501:PRO:HD3	1.87	0.56
1:B:142:ILE:HG12	1:B:170:GLU:HG2	1.87	0.56
1:B:383:ASN:HB3	1:B:624:GLN:O	2.05	0.56
1:C:822:LEU:HD11	1:C:825:CYS:HB2	1.86	0.56
1:C:355:ASN:OD1	1:C:388:ARG:HD3	2.06	0.55
1:D:794:GLY:HA3	5:D:4049:HOH:O	2.06	0.55
1:A:335:VAL:HG22	1:A:344:LEU:HD12	1.88	0.55
1:B:355:ASN:OD1	1:B:388:ARG:HD3	2.07	0.55
1:C:258:VAL:HG23	1:C:291:LEU:HD23	1.89	0.55
1:D:632:SER:HB2	5:D:4021:HOH:O	2.06	0.55
1:A:142:ILE:HG12	1:A:170:GLU:HG2	1.89	0.55
1:D:864:MET:HE3	1:D:866:ILE:HD11	1.88	0.55
1:C:38:ASN:HB3	1:C:41:GLU:HG3	1.89	0.55
1:C:58:TRP:CD1	1:C:86:VAL:HB	2.41	0.55
1:B:668:VAL:HG11	1:B:680:ILE:HG12	1.89	0.54
1:A:881:ARG:HE	1:A:987:ASP:CG	2.15	0.54
1:B:255:ARG:HB2	1:B:316:HIS:CE1	2.42	0.54
1:C:153:TRP:HB2	1:C:185:ALA:HB3	1.88	0.54
1:B:251:ARG:H	1:B:254:LEU:HD12	1.73	0.54
1:A:958:ASN:OD1	1:A:985:ASN:ND2	2.35	0.54
1:C:361:PRO:HD3	1:C:609:ALA:HB1	1.88	0.54
1:A:355:ASN:OD1	1:A:388:ARG:HD3	2.08	0.54
1:C:367:MET:HB3	1:C:372:MET:HE3	1.90	0.54
1:C:800:ARG:O	1:C:800:ARG:HG3	2.08	0.54
1:C:928:PRO:HB2	1:C:973:ARG:HH11	1.73	0.54
1:C:870:VAL:HG21	1:C:882:ILE:HG23	1.89	0.54
1:B:486:TYR:CE2	1:B:488:GLY:HA3	2.43	0.54
1:C:86:VAL:HG13	1:C:87:PRO:HA	1.90	0.53
1:D:333:ARG:HH11	1:D:333:ARG:HG2	1.73	0.53
1:B:750:GLU:HG2	1:B:755:ARG:HG3	1.90	0.53
1:C:19:PRO:HD3	1:C:112:PRO:HB3	1.91	0.53
1:C:262:GLN:OE1	1:C:299:LYS:HD2	2.08	0.53
1:D:646:HIS:HB3	5:D:4019:HOH:O	2.08	0.53
1:A:395:HIS:ND1	1:A:396:PRO:HD2	2.23	0.53
1:B:367:MET:HB3	1:B:372:MET:CE	2.38	0.53
1:D:837:THR:HG23	1:D:855:THR:HG22	1.90	0.53
1:B:942:ARG:HA	1:B:953:GLY:O	2.08	0.53
1:D:794:GLY:HA2	1:D:998:SER:O	2.09	0.52
1:B:768:MET:HB2	1:B:775:GLN:HB2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:613:PRO:HB3	1:A:617:LEU:CD2	2.40	0.52
1:B:427:THR:HG21	1:B:462:SER:HB3	1.91	0.52
1:D:942:ARG:HA	1:D:953:GLY:O	2.10	0.52
1:B:456:TRP:NE1	1:B:482:ARG:HD2	2.25	0.52
1:C:1000:SER:O	1:C:1001:PRO:C	2.53	0.51
1:D:99:ILE:O	1:D:204:ARG:HG2	2.11	0.51
1:D:373:VAL:O	1:D:377:LEU:HG	2.10	0.51
1:A:334:GLU:CG	4:A:8006:DMS:H13	2.40	0.51
1:B:226:HIS:CD2	1:B:448:ARG:HD3	2.45	0.51
1:C:38:ASN:HB3	1:C:41:GLU:CG	2.41	0.51
1:C:260:LEU:HD11	1:C:309:TYR:HB3	1.93	0.51
1:C:882:ILE:HD13	1:C:1014:TYR:CD1	2.46	0.51
1:B:879:PRO:O	1:B:1009:LEU:HD12	2.11	0.51
1:C:142:ILE:HG12	1:C:170:GLU:HG2	1.92	0.51
1:C:245:GLN:HG2	1:C:288:ARG:HG2	1.93	0.51
1:D:337:ILE:HA	1:D:341:LEU:O	2.11	0.51
1:C:330:VAL:HA	4:C:8001:DMS:H22	1.92	0.51
1:B:256:VAL:O	1:B:271:THR:HA	2.11	0.51
1:A:610:ASP:O	1:A:611:ARG:HB2	2.11	0.51
1:A:1020:TRP:HD1	1:A:1021:CYS:N	2.08	0.51
1:B:257:THR:HA	1:B:270:GLY:O	2.11	0.51
1:D:245:GLN:HG2	1:D:288:ARG:HG2	1.92	0.51
1:A:414:ASN:O	1:A:439:ARG:HD3	2.11	0.50
1:B:863:GLN:HG2	1:B:1021:CYS:HB3	1.93	0.50
1:A:146:VAL:HG22	1:A:208:ILE:HG12	1.92	0.50
1:A:854:LYS:HA	1:A:867:THR:O	2.12	0.50
1:B:607:VAL:HG12	1:B:613:PRO:HA	1.93	0.50
1:B:615:PRO:O	1:B:618:THR:HG22	2.10	0.50
1:D:980:GLU:HA	5:D:4079:HOH:O	2.10	0.50
1:B:372:MET:HE1	1:B:395:HIS:HB3	1.93	0.50
1:B:30:HIS:NE2	1:B:170:GLU:OE1	2.43	0.50
1:C:708:TRP:CE3	1:C:709:SER:HB3	2.46	0.50
1:A:763:GLY:HA3	1:A:822:LEU:HD13	1.94	0.50
1:B:143:PHE:O	1:B:168:PRO:HA	2.11	0.50
1:C:619:GLU:HA	1:C:912:ALA:HB2	1.92	0.50
1:C:991:MET:HE3	1:C:1003:VAL:HG21	1.94	0.50
1:D:784:PHE:HA	1:D:881:ARG:O	2.11	0.50
1:B:778:THR:HB	1:B:887:GLN:CB	2.42	0.50
1:A:341:LEU:HD13	1:A:348:PRO:HB3	1.94	0.49
1:D:708:TRP:CH2	4:D:8003:DMS:H23	2.47	0.49
1:D:279:ILE:HG12	1:D:284:GLY:HA2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:696:LEU:HB2	1:A:722:LEU:HD11	1.94	0.49
1:B:147:ASN:HB3	1:B:206:SER:HA	1.94	0.49
1:D:778:THR:H	1:D:887:GLN:HB3	1.77	0.49
1:B:50:GLN:O	1:B:215:LEU:HA	2.11	0.49
1:B:336:ARG:HB3	4:B:8006:DMS:H23	1.94	0.49
1:B:672:VAL:HG22	1:B:678:GLN:HB2	1.93	0.49
1:C:147:ASN:HB3	1:C:206:SER:HA	1.94	0.49
1:D:133:TRP:HA	1:D:216:HIS:CE1	2.47	0.49
1:D:536:CYS:O	1:D:537:GLU:HG3	2.12	0.49
1:D:537:GLU:HA	1:D:566:PHE:O	2.12	0.49
1:B:17:GLU:HG2	1:B:114:VAL:HG22	1.95	0.48
1:D:237:ARG:HG2	1:D:296:GLU:OE1	2.13	0.48
1:A:878:HIS:CE1	1:A:1010:SER:HB3	2.47	0.48
1:C:124:SER:HA	1:C:184:LEU:O	2.13	0.48
1:C:606:LEU:O	1:C:614:HIS:HB2	2.13	0.48
1:A:333:ARG:HA	1:A:345:ASN:OD1	2.14	0.48
1:A:343:LEU:HA	1:A:347:LYS:O	2.13	0.48
1:D:333:ARG:NH1	1:D:333:ARG:HG2	2.28	0.48
1:C:88:SER:HA	1:C:366:VAL:HG21	1.95	0.48
1:C:163:GLN:O	1:C:164:ASP:HB3	2.13	0.48
1:B:377:LEU:HD22	1:B:708:TRP:HA	1.96	0.48
1:D:598:ASP:O	1:D:601:PHE:HB2	2.13	0.48
1:A:114:VAL:HA	5:A:4077:HOH:O	2.13	0.48
1:D:573:GLN:HB2	1:D:602:CYS:O	2.13	0.48
1:C:897:TRP:CZ2	1:C:938:ARG:HG2	2.49	0.48
1:B:763:GLY:HA3	1:B:822:LEU:HD13	1.96	0.48
1:A:524:LEU:HD11	1:A:562:LEU:HG	1.96	0.47
1:A:907:PRO:HG2	1:A:990:HIS:O	2.14	0.47
1:A:147:ASN:ND2	1:A:205:MET:O	2.46	0.47
1:C:496:THR:OG1	1:C:498:ILE:O	2.31	0.47
1:D:132:SER:HA	1:D:135:GLN:HB2	1.96	0.47
1:A:146:VAL:HG11	1:A:150:PHE:CD2	2.49	0.47
1:D:842:TRP:HZ3	1:D:852:SER:HB3	1.79	0.47
1:B:875:ASP:OD2	1:B:875:ASP:N	2.43	0.47
1:C:440:VAL:HG13	1:C:475:ILE:HD11	1.96	0.47
1:D:226:HIS:O	1:D:242:ALA:HA	2.14	0.47
1:A:140:ARG:HG2	1:A:215:LEU:HB3	1.97	0.47
1:C:47:PRO:HA	4:C:8011:DMS:H22	1.96	0.47
1:C:738:PRO:HG3	1:C:751:LEU:HD13	1.97	0.47
1:D:603:MET:HE1	1:D:930:VAL:HB	1.97	0.47
1:B:753:ASN:HA	5:B:4072:HOH:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:824:GLN:HB3	1:C:839:ALA:HB3	1.96	0.47
1:A:245:GLN:HG2	1:A:288:ARG:CG	2.41	0.47
1:A:697:THR:HG23	1:A:719:GLN:HG3	1.96	0.47
1:A:1000:SER:O	1:A:1001:PRO:C	2.57	0.47
1:C:786:ARG:NH2	1:C:991:MET:HE2	2.30	0.47
1:D:347:LYS:HB3	1:D:643:LEU:HD22	1.97	0.46
1:C:906:TYR:HB3	1:C:907:PRO:HD2	1.96	0.46
1:B:390:SER:HA	1:B:391:HIS:HA	1.70	0.46
1:B:702:GLN:O	1:B:712:GLY:N	2.44	0.46
1:C:942:ARG:O	1:C:952:ARG:HA	2.16	0.46
1:C:598:ASP:O	1:C:601:PHE:HB2	2.16	0.46
1:D:105:TYR:CE2	1:D:199:ASP:HB2	2.51	0.46
1:D:906:TYR:OH	1:D:934:GLU:HG3	2.16	0.46
1:B:610:ASP:O	1:B:611:ARG:HB2	2.16	0.46
1:D:153:TRP:HB2	1:D:185:ALA:HB3	1.97	0.46
1:D:30:HIS:ND1	1:D:31:PRO:O	2.46	0.46
1:B:344:LEU:HD23	1:B:644:PHE:CZ	2.51	0.46
1:C:873:ALA:O	1:C:876:THR:HG22	2.16	0.46
1:A:227:VAL:HG13	1:A:240:LEU:HD21	1.98	0.46
1:B:957:PHE:CD1	1:B:957:PHE:C	2.94	0.46
1:D:163:GLN:O	1:D:164:ASP:HB3	2.16	0.46
1:D:220:THR:HG21	1:D:254:LEU:HD11	1.97	0.46
1:A:107:ILE:HB	5:A:4077:HOH:O	2.14	0.46
1:A:486:TYR:CZ	1:A:488:GLY:HA3	2.51	0.46
1:B:91:GLN:HG3	1:B:96:ASP:OD1	2.16	0.46
1:B:150:PHE:HA	1:B:187:MET:O	2.16	0.45
1:D:35:SER:O	1:D:50:GLN:HG3	2.16	0.45
1:C:728:VAL:HG12	1:D:851:ILE:HD11	1.98	0.45
1:D:708:TRP:CE3	1:D:709:SER:HB3	2.51	0.45
1:B:749:ILE:HD11	1:B:836:ILE:HD11	1.97	0.45
1:B:861:SER:HB3	5:B:4061:HOH:O	2.15	0.45
1:B:896:ASN:HB3	1:B:945:ASN:HB2	1.97	0.45
1:D:348:PRO:O	1:D:643:LEU:HD11	2.16	0.45
1:A:118:ASN:O	1:A:119:PRO:C	2.60	0.45
1:A:895:VAL:HG11	1:A:925:MET:HE2	1.97	0.45
1:C:492:ASP:HB3	1:C:499:ILE:HG23	1.98	0.45
1:D:892:ALA:HB3	1:D:946:TYR:CZ	2.50	0.45
1:B:928:PRO:HB2	1:B:973:ARG:HH11	1.82	0.45
1:D:338:GLU:HB2	5:D:4031:HOH:O	2.15	0.45
1:B:755:ARG:HB3	1:B:769:TRP:HB2	1.99	0.45
1:A:240:LEU:HD23	1:A:241:GLU:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:MET:C	1:B:246:MET:SD	3.00	0.45
1:A:390:SER:HB2	1:A:391:HIS:CE1	2.52	0.45
1:A:390:SER:HA	1:A:391:HIS:HA	1.74	0.45
1:C:60:PHE:HA	1:C:122:CYS:O	2.17	0.45
1:C:367:MET:HB3	1:C:372:MET:CE	2.47	0.45
1:D:43:ARG:HD2	1:D:261:TRP:CD2	2.52	0.45
1:B:58:TRP:CD1	1:B:86:VAL:HB	2.52	0.45
1:C:644:PHE:O	1:C:674:PRO:HG3	2.17	0.45
1:C:881:ARG:HE	1:C:987:ASP:CG	2.25	0.44
1:C:883:GLY:HA3	1:C:987:ASP:HA	1.99	0.44
1:D:88:SER:HA	1:D:366:VAL:CG2	2.45	0.44
1:D:410:VAL:HG22	1:D:455:ILE:HB	1.99	0.44
1:A:127:PHE:HE2	1:A:184:LEU:HG	1.82	0.44
1:C:615:PRO:HB3	1:C:904:GLU:HB3	1.99	0.44
1:A:613:PRO:HB3	1:A:617:LEU:HD22	2.00	0.44
1:B:788:PRO:HD2	1:B:968:MET:HB2	1.98	0.44
1:B:456:TRP:CE2	1:B:482:ARG:HD2	2.53	0.44
1:B:766:SER:HA	1:B:779:PRO:HB3	1.99	0.44
1:D:335:VAL:HG22	1:D:344:LEU:HD12	1.99	0.44
1:D:698:VAL:O	1:D:717:TRP:HA	2.16	0.44
1:B:870:VAL:HG21	1:B:882:ILE:CG2	2.47	0.44
1:D:545:SER:O	1:D:909:ARG:HD3	2.18	0.44
1:D:789:LEU:HD21	1:D:934:GLU:HB3	2.00	0.44
1:A:127:PHE:CE2	1:A:184:LEU:HG	2.53	0.43
1:A:770:ILE:HD13	1:A:1022:GLN:HG3	1.99	0.43
1:C:105:TYR:CE2	1:C:199:ASP:HB2	2.53	0.43
1:D:217:LYS:NZ	1:D:326:GLU:OE2	2.50	0.43
1:A:600:GLN:HB2	1:A:603:MET:HE2	2.01	0.43
1:B:737:ILE:HD12	1:B:738:PRO:HD2	2.00	0.43
1:A:753:ASN:HB2	1:A:771:GLY:HA2	1.99	0.43
1:B:549:PHE:CE2	1:B:620:ALA:HA	2.53	0.43
1:C:141:ILE:HB	1:C:173:LEU:HD11	1.99	0.43
1:C:553:TRP:CZ2	1:C:624:GLN:HG2	2.52	0.43
1:C:963:SER:HB3	1:C:983:TRP:CE2	2.53	0.43
1:D:693:GLN:HB2	1:D:724:GLU:HG3	2.01	0.43
1:A:256:VAL:HG23	1:A:274:PHE:CZ	2.53	0.43
1:B:361:PRO:HB2	1:B:576:ILE:HG12	2.00	0.43
1:B:429:ASP:OD1	1:B:431:ARG:HG3	2.19	0.43
1:B:778:THR:HB	1:B:887:GLN:HB3	2.01	0.43
1:C:854:LYS:HG2	1:C:868:VAL:HG22	2.00	0.43
1:D:693:GLN:HG2	1:D:695:TRP:NE1	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:ARG:HH12	1:A:264:GLU:HG2	1.83	0.43
1:A:354:VAL:O	1:A:387:VAL:HG23	2.19	0.43
1:A:613:PRO:CB	1:A:617:LEU:HD23	2.49	0.43
1:B:390:SER:HB2	1:B:391:HIS:CE1	2.53	0.43
1:A:738:PRO:HG3	1:A:751:LEU:HD13	1.99	0.43
1:B:794:GLY:HA2	1:B:998:SER:O	2.19	0.43
1:B:1000:SER:O	1:B:1001:PRO:C	2.62	0.43
1:C:522:LYS:HE3	1:D:560:PRO:HD3	2.00	0.43
1:D:778:THR:HA	1:D:779:PRO:HD3	1.85	0.43
1:B:474:TRP:CE2	1:B:478:VAL:HG21	2.54	0.43
1:D:544:ASN:HB3	1:D:789:LEU:HD13	2.01	0.43
1:D:549:PHE:CE2	1:D:620:ALA:HA	2.54	0.43
1:D:608:PHE:CZ	1:D:614:HIS:HD2	2.37	0.43
1:A:260:LEU:HD11	1:A:309:TYR:HB3	2.00	0.43
1:B:443:MET:HE3	1:B:456:TRP:CE3	2.54	0.43
1:A:147:ASN:HB3	1:A:206:SER:HA	2.00	0.42
1:A:454:ILE:HG13	1:A:455:ILE:HG13	2.01	0.42
1:B:505:ARG:O	1:B:519:SER:HA	2.19	0.42
1:C:414:ASN:C	1:C:415:ILE:HG13	2.45	0.42
1:D:225:PHE:HA	1:D:243:GLU:O	2.19	0.42
1:A:59:ARG:HB2	1:A:124:SER:OG	2.19	0.42
1:A:541:ALA:HB2	1:A:546:LEU:HD13	2.01	0.42
1:C:961:ARG:NH2	1:C:979:GLU:O	2.52	0.42
1:B:496:THR:OG1	1:B:498:ILE:O	2.35	0.42
1:B:906:TYR:HB3	1:B:907:PRO:HD2	2.02	0.42
1:C:755:ARG:HB3	1:C:769:TRP:HE3	1.84	0.42
1:D:240:LEU:HD13	1:D:260:LEU:HD22	2.01	0.42
1:A:226:HIS:CD2	1:A:448:ARG:HD3	2.55	0.42
1:B:235:PHE:HB3	1:B:300:LEU:HD11	2.01	0.42
1:D:292:ARG:N	4:D:8004:DMS:O	2.50	0.42
1:D:500:CYS:HA	1:D:534:ILE:O	2.20	0.42
1:D:595:THR:HA	1:D:596:PRO:C	2.45	0.42
1:D:927:THR:HA	1:D:928:PRO:HD3	1.91	0.42
1:D:972:HIS:HB2	1:D:975:LEU:HD12	2.01	0.42
1:B:249:GLU:OE2	1:B:251:ARG:HD3	2.19	0.42
1:B:890:GLN:HE21	1:B:891:VAL:C	2.28	0.42
1:C:350:LEU:HD11	1:C:556:PHE:HB3	2.01	0.42
1:C:379:MET:HE1	1:C:387:VAL:HB	2.01	0.42
1:C:896:ASN:HA	1:C:918:TRP:O	2.20	0.42
1:C:939:CYS:HA	1:C:956:GLN:HB3	2.02	0.42
1:B:524:LEU:HD11	1:B:562:LEU:HG	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:937:LEU:HA	1:B:957:PHE:O	2.19	0.42
1:C:525:SER:O	1:D:561:ARG:HD3	2.20	0.42
1:D:23:GLN:C	1:D:23:GLN:CD	2.88	0.42
1:D:272:ALA:HB1	1:D:273:PRO:HD2	2.02	0.42
1:D:569:ASP:O	1:D:605:GLY:HA2	2.20	0.42
1:A:931:PHE:HA	1:A:932:PRO:HD3	1.86	0.42
1:C:603:MET:HE1	1:C:930:VAL:HB	2.00	0.42
1:C:864:MET:HE3	1:C:866:ILE:HD11	2.02	0.42
1:D:379:MET:HE1	1:D:387:VAL:HB	2.01	0.42
1:A:43:ARG:HD2	1:A:261:TRP:CD2	2.55	0.41
1:A:429:ASP:HA	1:A:430:PRO:HD2	1.89	0.41
1:B:545:SER:OG	1:B:546:LEU:N	2.52	0.41
1:B:957:PHE:HA	1:B:985:ASN:O	2.20	0.41
1:D:458:LEU:HD11	1:D:472:TYR:HB2	2.02	0.41
1:A:807:VAL:HG13	1:A:808:GLU:HG2	2.01	0.41
1:B:50:GLN:N	1:B:50:GLN:CD	2.78	0.41
1:A:855:THR:OG1	1:A:867:THR:HB	2.20	0.41
1:B:147:ASN:HA	1:B:148:SER:HA	1.76	0.41
1:B:559:TYR:HB2	1:B:562:LEU:HD12	2.02	0.41
1:C:414:ASN:O	1:C:439:ARG:HD3	2.20	0.41
1:C:997:ASP:HB2	1:C:999:TRP:CE2	2.55	0.41
1:A:655:MET:HE1	1:A:662:PRO:HG3	2.02	0.41
1:A:824:GLN:O	1:A:838:THR:HA	2.20	0.41
1:C:142:ILE:HG21	5:C:4015:HOH:O	2.19	0.41
1:C:411:ASP:HB2	1:C:453:VAL:HG13	2.02	0.41
1:B:274:PHE:HA	1:B:289:VAL:HB	2.03	0.41
1:B:613:PRO:HB3	1:B:617:LEU:HD23	2.02	0.41
1:B:759:ASN:ND2	5:B:4104:HOH:O	2.54	0.41
1:C:147:ASN:OD1	1:C:148:SER:HB3	2.21	0.41
1:D:354:VAL:CG1	1:D:379:MET:HE3	2.50	0.41
1:C:251:ARG:H	1:C:254:LEU:HD12	1.85	0.41
1:C:421:VAL:HA	1:C:422:PRO:HA	1.94	0.41
1:D:218:PRO:HB2	1:D:220:THR:O	2.21	0.41
1:C:796:SER:HB2	1:C:802:ASP:HB3	2.03	0.41
1:A:33:PHE:HB3	1:A:326:GLU:OE2	2.20	0.41
1:A:58:TRP:O	1:A:83:THR:HA	2.21	0.41
1:A:523:TRP:CD1	1:A:526:LEU:HD12	2.55	0.41
1:A:526:LEU:O	1:A:527:PRO:C	2.64	0.41
1:A:957:PHE:HA	1:A:985:ASN:O	2.21	0.41
1:A:1020:TRP:CD1	1:A:1020:TRP:C	2.98	0.41
1:C:147:ASN:HA	1:C:148:SER:HA	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:984:LEU:HD21	1:C:986:ILE:HD11	2.02	0.41
1:D:257:THR:HA	1:D:270:GLY:O	2.21	0.41
1:D:390:SER:HA	1:D:391:HIS:HA	1.76	0.41
1:D:615:PRO:O	1:D:618:THR:HG22	2.21	0.41
1:A:568:TRP:HA	1:A:569:ASP:HA	1.81	0.41
1:A:737:ILE:HA	1:A:738:PRO:HD3	1.87	0.41
1:B:18:ASN:HA	1:B:19:PRO:HD2	1.93	0.41
1:A:542:MET:HA	1:A:604:ASN:HA	2.03	0.40
1:B:835:LEU:HA	1:B:856:TYR:O	2.21	0.40
1:D:89:ASN:HA	1:D:206:SER:O	2.21	0.40
1:B:376:ILE:CD1	1:B:401:LEU:HB3	2.51	0.40
1:C:46:ARG:HG3	1:C:47:PRO:HD2	2.03	0.40
1:C:190:ARG:HG2	1:C:191:TRP:CE2	2.56	0.40
1:C:260:LEU:HD11	1:C:309:TYR:CB	2.51	0.40
1:D:92:MET:HA	1:D:92:MET:HE2	2.03	0.40
1:D:262:GLN:NE2	1:D:299:LYS:HD2	2.36	0.40
1:A:107:ILE:HD12	5:A:4077:HOH:O	2.19	0.40
1:A:380:LYS:O	4:A:8003:DMS:H11	2.21	0.40
1:B:835:LEU:HD11	1:B:855:THR:HB	2.02	0.40
1:D:577:LYS:O	1:D:584:PRO:HA	2.21	0.40
1:D:696:LEU:O	1:D:719:GLN:HA	2.21	0.40
1:B:568:TRP:CG	1:B:569:ASP:HB3	2.56	0.40
1:C:380:LYS:NZ	1:C:406:GLY:O	2.55	0.40
1:C:780:LEU:HA	1:C:886:CYS:HB3	2.04	0.40
1:D:147:ASN:O	1:D:206:SER:HB2	2.21	0.40
1:A:421:VAL:HA	1:A:422:PRO:HA	1.97	0.40
1:C:607:VAL:HG12	1:C:613:PRO:HA	2.04	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	1009/1052 (96%)	949 (94%)	58 (6%)	2 (0%)	43	72
1	B	1009/1052 (96%)	948 (94%)	57 (6%)	4 (0%)	30	60
1	C	1009/1052 (96%)	953 (94%)	52 (5%)	4 (0%)	30	60
1	D	1009/1052 (96%)	943 (94%)	65 (6%)	1 (0%)	48	77
All	All	4036/4208 (96%)	3793 (94%)	232 (6%)	11 (0%)	36	66

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	164	ASP
1	B	461	GLU
1	A	541	ALA
1	C	733	ALA
1	B	540	HIS
1	C	164	ASP
1	D	109	VAL
1	A	539	ALA
1	C	539	ALA
1	B	633	GLY
1	C	1001	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	864/898 (96%)	840 (97%)	24 (3%)	38	73
1	B	864/898 (96%)	843 (98%)	21 (2%)	43	77
1	C	864/898 (96%)	837 (97%)	27 (3%)	35	70
1	D	864/898 (96%)	842 (98%)	22 (2%)	42	76
All	All	3456/3592 (96%)	3362 (97%)	94 (3%)	39	74

All (94) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	80	GLU
1	A	131	GLU
1	A	195	SER
1	A	213	SER
1	A	291	LEU
1	A	293	LEU
1	A	515	VAL
1	A	519	SER
1	A	595	THR
1	A	604	ASN
1	A	655	MET
1	A	663	LEU
1	A	681	GLU
1	A	730	LEU
1	A	737	ILE
1	A	745	MET
1	A	773	LYS
1	A	817	GLN
1	A	855	THR
1	A	910	LEU
1	A	916	ASP
1	A	924	ASP
1	A	956	GLN
1	B	13	ARG
1	B	80	GLU
1	B	126	THR
1	B	186	VAL
1	B	267	VAL
1	B	269	SER
1	B	344	LEU
1	B	519	SER
1	B	535	LEU
1	B	546	LEU
1	B	635	THR
1	B	663	LEU
1	B	743	SER
1	B	748	CYS
1	B	774	LYS
1	B	819	GLU
1	B	829	THR
1	B	876	THR
1	B	965	GLN

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Mol	Chain	Res	Type
1	B	968	MET
1	B	1004	SER
1	C	46	ARG
1	C	71	GLU
1	C	80	GLU
1	C	126	THR
1	C	129	VAL
1	C	181	GLU
1	C	213	SER
1	C	219	THR
1	C	291	LEU
1	C	546	LEU
1	C	635	THR
1	C	663	LEU
1	C	672	VAL
1	C	729	THR
1	C	737	ILE
1	C	741	THR
1	C	751	LEU
1	C	773	LYS
1	C	785	THR
1	C	789	LEU
1	C	800	ARG
1	C	819	GLU
1	C	890	GLN
1	C	937	LEU
1	C	956	GLN
1	C	1002	SER
1	C	1022	GLN
1	D	80	GLU
1	D	129	VAL
1	D	135	GLN
1	D	213	SER
1	D	293	LEU
1	D	333	ARG
1	D	519	SER
1	D	546	LEU
1	D	580	GLU
1	D	612	THR
1	D	652	LEU
1	D	663	LEU
1	D	672	VAL

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Mol	Chain	Res	Type
1	D	728	VAL
1	D	744	GLU
1	D	748	CYS
1	D	799	THR
1	D	824	GLN
1	D	956	GLN
1	D	966	GLN
1	D	970	THR
1	D	986	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	128	ASN
1	A	163	GLN
1	A	394	ASN
1	A	581	ASN
1	A	614	HIS
1	A	634	GLN
1	A	704	ASN
1	A	824	GLN
1	A	949	HIS
1	A	977	HIS
1	B	50	GLN
1	B	374	GLN
1	B	391	HIS
1	B	394	ASN
1	B	485	GLN
1	B	624	GLN
1	B	702	GLN
1	B	844	HIS
1	B	890	GLN
1	B	949	HIS
1	B	956	GLN
1	B	977	HIS
1	C	128	ASN
1	C	155	ASN
1	C	381	GLN
1	C	485	GLN
1	C	704	ASN
1	C	804	ASN

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Mol	Chain	Res	Type
1	C	817	GLN
1	C	885	ASN
1	C	956	GLN
1	C	977	HIS
1	C	1022	GLN
1	D	93	HIS
1	D	363	HIS
1	D	382	ASN
1	D	464	HIS
1	D	485	GLN
1	D	510	GLN
1	D	614	HIS
1	D	622	HIS
1	D	628	GLN
1	D	739	HIS
1	D	804	ASN
1	D	824	GLN
1	D	844	HIS
1	D	903	GLN
1	D	945	ASN
1	D	950	GLN
1	D	956	GLN
1	D	974	HIS
1	D	1015	HIS

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 64 ligands modelled in this entry, 19 are monoatomic - leaving 45 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	A	8002	-	3,3,3	2.72	1 (33%)	3,3,3	0.74	0
4	DMS	D	8008	-	3,3,3	2.71	1 (33%)	3,3,3	0.54	0
4	DMS	B	8004	-	3,3,3	2.80	1 (33%)	3,3,3	0.61	0
4	DMS	C	8012	-	3,3,3	2.77	1 (33%)	3,3,3	0.59	0
4	DMS	B	8008	-	3,3,3	2.71	1 (33%)	3,3,3	0.58	0
4	DMS	A	8007	-	3,3,3	2.80	1 (33%)	3,3,3	0.68	0
4	DMS	C	8002	-	3,3,3	2.81	1 (33%)	3,3,3	0.89	0
4	DMS	A	8008	-	3,3,3	2.76	1 (33%)	3,3,3	0.58	0
4	DMS	A	8010	-	3,3,3	2.74	1 (33%)	3,3,3	0.67	0
4	DMS	D	8003	-	3,3,3	2.67	1 (33%)	3,3,3	0.57	0
4	DMS	C	8003	-	3,3,3	2.73	1 (33%)	3,3,3	0.61	0
4	DMS	D	8001	-	3,3,3	2.74	1 (33%)	3,3,3	0.64	0
4	DMS	C	8008	-	3,3,3	2.66	1 (33%)	3,3,3	0.63	0
4	DMS	D	8010	-	3,3,3	2.75	1 (33%)	3,3,3	0.67	0
4	DMS	C	8011	-	3,3,3	2.72	1 (33%)	3,3,3	0.47	0
4	DMS	A	8003	-	3,3,3	2.72	1 (33%)	3,3,3	0.57	0
4	DMS	B	8009	-	3,3,3	2.75	1 (33%)	3,3,3	0.61	0
4	DMS	A	8005	-	3,3,3	2.75	1 (33%)	3,3,3	0.69	0
4	DMS	C	8009	-	3,3,3	2.76	1 (33%)	3,3,3	0.62	0
4	DMS	D	8009	-	3,3,3	2.67	1 (33%)	3,3,3	0.52	0
4	DMS	C	8004	-	3,3,3	2.66	1 (33%)	3,3,3	0.49	0
4	DMS	D	8004	-	3,3,3	2.63	1 (33%)	3,3,3	0.71	0
4	DMS	A	8012	-	3,3,3	2.77	1 (33%)	3,3,3	0.62	0
4	DMS	A	8006	-	3,3,3	2.75	1 (33%)	3,3,3	0.61	0
4	DMS	D	8002	-	3,3,3	2.68	1 (33%)	3,3,3	0.55	0
4	DMS	D	8007	-	3,3,3	2.80	1 (33%)	3,3,3	0.83	0
4	DMS	B	8006	-	3,3,3	2.74	1 (33%)	3,3,3	0.66	0
4	DMS	C	8005	-	3,3,3	2.79	1 (33%)	3,3,3	0.65	0
4	DMS	C	8006	-	3,3,3	2.74	1 (33%)	3,3,3	0.60	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DMS	B	8007	-	3,3,3	2.72	1 (33%)	3,3,3	0.49	0
4	DMS	B	8010	-	3,3,3	2.71	1 (33%)	3,3,3	0.49	0
4	DMS	A	8009	-	3,3,3	2.73	1 (33%)	3,3,3	0.52	0
4	DMS	B	8003	-	3,3,3	2.75	1 (33%)	3,3,3	0.61	0
4	DMS	D	8006	-	3,3,3	2.77	1 (33%)	3,3,3	0.64	0
4	DMS	B	8001	-	3,3,3	2.70	1 (33%)	3,3,3	0.63	0
4	DMS	C	8010	-	3,3,3	2.73	1 (33%)	3,3,3	0.82	0
4	DMS	C	8001	-	3,3,3	2.60	1 (33%)	3,3,3	0.47	0
4	DMS	D	8005	-	3,3,3	2.74	1 (33%)	3,3,3	0.64	0
4	DMS	B	8002	-	3,3,3	2.72	1 (33%)	3,3,3	0.59	0
4	DMS	A	8004	-	3,3,3	2.68	1 (33%)	3,3,3	0.50	0
4	DMS	B	8011	-	3,3,3	2.72	1 (33%)	3,3,3	0.63	0
4	DMS	C	8007	-	3,3,3	2.78	1 (33%)	3,3,3	0.62	0
4	DMS	B	8005	-	3,3,3	2.71	1 (33%)	3,3,3	0.51	0
4	DMS	A	8011	-	3,3,3	2.77	1 (33%)	3,3,3	0.58	0
4	DMS	A	8001	-	3,3,3	2.69	1 (33%)	3,3,3	0.55	0

All (45) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	8002	DMS	O-S	4.72	1.81	1.50
4	B	8004	DMS	O-S	4.70	1.81	1.50
4	D	8007	DMS	O-S	4.68	1.81	1.50
4	C	8005	DMS	O-S	4.68	1.81	1.50
4	A	8007	DMS	O-S	4.68	1.81	1.50
4	D	8006	DMS	O-S	4.66	1.81	1.50
4	C	8007	DMS	O-S	4.65	1.80	1.50
4	A	8011	DMS	O-S	4.65	1.80	1.50
4	C	8012	DMS	O-S	4.64	1.80	1.50
4	A	8006	DMS	O-S	4.63	1.80	1.50
4	A	8008	DMS	O-S	4.63	1.80	1.50
4	A	8005	DMS	O-S	4.62	1.80	1.50
4	A	8012	DMS	O-S	4.62	1.80	1.50
4	C	8009	DMS	O-S	4.62	1.80	1.50
4	D	8001	DMS	O-S	4.62	1.80	1.50
4	B	8009	DMS	O-S	4.60	1.80	1.50
4	B	8003	DMS	O-S	4.60	1.80	1.50
4	C	8006	DMS	O-S	4.59	1.80	1.50
4	B	8006	DMS	O-S	4.59	1.80	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	8010	DMS	O-S	4.59	1.80	1.50
4	C	8003	DMS	O-S	4.58	1.80	1.50
4	D	8010	DMS	O-S	4.58	1.80	1.50
4	A	8009	DMS	O-S	4.58	1.80	1.50
4	D	8005	DMS	O-S	4.57	1.80	1.50
4	C	8010	DMS	O-S	4.57	1.80	1.50
4	B	8011	DMS	O-S	4.57	1.80	1.50
4	A	8002	DMS	O-S	4.57	1.80	1.50
4	A	8003	DMS	O-S	4.56	1.80	1.50
4	C	8011	DMS	O-S	4.56	1.80	1.50
4	B	8007	DMS	O-S	4.55	1.80	1.50
4	B	8010	DMS	O-S	4.55	1.80	1.50
4	B	8002	DMS	O-S	4.54	1.80	1.50
4	B	8005	DMS	O-S	4.54	1.80	1.50
4	D	8008	DMS	O-S	4.54	1.80	1.50
4	B	8008	DMS	O-S	4.54	1.80	1.50
4	B	8001	DMS	O-S	4.52	1.80	1.50
4	A	8001	DMS	O-S	4.51	1.80	1.50
4	D	8002	DMS	O-S	4.51	1.80	1.50
4	A	8004	DMS	O-S	4.50	1.79	1.50
4	D	8009	DMS	O-S	4.49	1.79	1.50
4	D	8003	DMS	O-S	4.48	1.79	1.50
4	C	8008	DMS	O-S	4.44	1.79	1.50
4	C	8004	DMS	O-S	4.43	1.79	1.50
4	D	8004	DMS	O-S	4.38	1.79	1.50
4	C	8001	DMS	O-S	4.36	1.79	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	8002	DMS	1	0
4	D	8003	DMS	1	0
4	D	8001	DMS	1	0
4	C	8011	DMS	1	0
4	A	8003	DMS	1	0
4	D	8004	DMS	1	0
4	A	8006	DMS	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	8006	DMS	1	0
4	C	8001	DMS	1	0
4	B	8011	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	1011/1052 (96%)	-0.11	12 (1%) 76 68	8, 25, 52, 91	0
1	B	1011/1052 (96%)	0.05	19 (1%) 66 57	9, 31, 52, 90	0
1	C	1011/1052 (96%)	0.11	32 (3%) 50 40	13, 30, 54, 92	0
1	D	1011/1052 (96%)	0.09	27 (2%) 56 46	6, 28, 61, 96	0
All	All	4044/4208 (96%)	0.03	90 (2%) 62 52	6, 28, 55, 96	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	688	PRO	4.6
1	C	736	ALA	4.5
1	A	733	ALA	3.7
1	B	733	ALA	3.6
1	D	766	SER	3.5
1	B	685	LEU	3.5
1	B	858	ILE	3.5
1	C	986	ILE	3.5
1	D	685	LEU	3.5
1	C	632	SER	3.4
1	A	581	ASN	3.4
1	D	686	PRO	3.4
1	C	843	GLN	3.3
1	C	745	MET	3.2
1	C	735	HIS	3.2
1	D	780	LEU	3.1
1	C	733	ALA	3.1
1	B	730	LEU	3.1
1	B	686	PRO	3.1
1	D	770	ILE	3.1
1	A	548	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	744	GLU	3.0
1	D	742	THR	3.0
1	A	735	HIS	3.0
1	B	32	PRO	3.0
1	D	769	TRP	2.9
1	D	731	PRO	2.9
1	A	462	SER	2.8
1	B	688	PRO	2.8
1	D	819	GLU	2.7
1	B	734	SER	2.7
1	A	736	ALA	2.7
1	A	737	ILE	2.7
1	D	737	ILE	2.7
1	C	128	ASN	2.7
1	C	686	PRO	2.7
1	C	863	GLN	2.7
1	C	171	PHE	2.6
1	B	148	SER	2.6
1	C	148	SER	2.6
1	D	768	MET	2.6
1	D	858	ILE	2.6
1	B	128	ASN	2.6
1	D	818	ALA	2.6
1	B	729	THR	2.6
1	C	716	ALA	2.5
1	D	733	ALA	2.5
1	C	688	PRO	2.5
1	A	181	GLU	2.5
1	D	767	GLN	2.5
1	C	78	LEU	2.5
1	C	1019	VAL	2.4
1	D	738	PRO	2.4
1	B	181	GLU	2.4
1	C	730	LEU	2.4
1	B	834	VAL	2.4
1	C	22	THR	2.3
1	D	740	LEU	2.3
1	C	662	PRO	2.3
1	C	731	PRO	2.3
1	B	728	VAL	2.3
1	D	933	SER	2.3
1	C	800	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	764	PHE	2.3
1	C	732	ALA	2.3
1	D	782	ASP	2.2
1	C	699	ARG	2.2
1	C	664	ALA	2.2
1	C	320	GLY	2.2
1	B	732	ALA	2.2
1	C	180	GLY	2.2
1	D	70	PRO	2.2
1	D	834	VAL	2.2
1	A	685	LEU	2.2
1	B	731	PRO	2.1
1	C	847	LYS	2.1
1	C	846	GLY	2.1
1	A	683	PRO	2.1
1	A	731	PRO	2.1
1	C	728	VAL	2.1
1	D	764	PHE	2.1
1	D	743	SER	2.1
1	D	109	VAL	2.1
1	C	931	PHE	2.0
1	C	691	ALA	2.0
1	D	891	VAL	2.0
1	A	741	THR	2.0
1	B	131	GLU	2.0
1	B	859	ASP	2.0
1	C	329	ASP	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($\text{\AA}^2$)	Q<0.9
4	DMS	A	8007	4/4	0.45	0.38	82,82,83,87	0
4	DMS	D	8010	4/4	0.68	0.34	68,69,70,72	0
4	DMS	B	8004	4/4	0.73	0.28	67,67,67,68	0
4	DMS	C	8006	4/4	0.83	0.23	82,83,83,83	0
4	DMS	D	8007	4/4	0.83	0.17	39,40,45,46	0
3	NA	B	3104	1/1	0.83	0.15	38,38,38,38	0
4	DMS	B	8005	4/4	0.87	0.23	83,83,83,83	0
4	DMS	D	8005	4/4	0.87	0.16	49,51,52,53	0
4	DMS	B	8011	4/4	0.88	0.18	70,71,72,72	0
4	DMS	A	8011	4/4	0.88	0.24	56,57,57,61	0
4	DMS	A	8005	4/4	0.88	0.22	62,63,64,66	0
2	MG	B	3001	1/1	0.88	0.18	45,45,45,45	0
4	DMS	B	8008	4/4	0.88	0.18	54,55,55,56	0
4	DMS	C	8003	4/4	0.89	0.22	77,77,77,77	0
4	DMS	A	8012	4/4	0.89	0.18	56,56,56,57	0
4	DMS	B	8003	4/4	0.89	0.17	46,46,48,48	0
4	DMS	B	8009	4/4	0.89	0.21	63,65,65,66	0
3	NA	B	3103	1/1	0.89	0.14	41,41,41,41	0
4	DMS	A	8008	4/4	0.90	0.18	62,64,64,65	0
4	DMS	C	8012	4/4	0.90	0.25	65,66,66,66	0
4	DMS	B	8007	4/4	0.90	0.16	53,55,55,56	0
4	DMS	A	8010	4/4	0.90	0.17	56,59,60,61	0
4	DMS	C	8005	4/4	0.90	0.14	54,54,56,58	0
4	DMS	A	8006	4/4	0.91	0.16	40,44,46,51	0
4	DMS	C	8010	4/4	0.91	0.20	53,54,54,58	0
4	DMS	A	8009	4/4	0.91	0.16	62,63,64,65	0
3	NA	C	3105	1/1	0.92	0.13	38,38,38,38	0
3	NA	D	3101	1/1	0.92	0.07	36,36,36,36	0
4	DMS	C	8007	4/4	0.92	0.12	43,43,44,48	0
2	MG	A	3001	1/1	0.92	0.15	42,42,42,42	0
3	NA	A	3101	1/1	0.93	0.06	26,26,26,26	0
2	MG	C	3001	1/1	0.93	0.14	43,43,43,43	0
4	DMS	C	8002	4/4	0.93	0.14	37,38,38,39	0
4	DMS	B	8006	4/4	0.93	0.13	52,52,52,55	0
4	DMS	A	8003	4/4	0.93	0.13	32,32,33,34	0
2	MG	C	3002	1/1	0.93	0.05	29,29,29,29	0
2	MG	B	3002	1/1	0.94	0.05	33,33,33,33	0
4	DMS	C	8011	4/4	0.94	0.11	46,47,47,48	0
4	DMS	B	8001	4/4	0.94	0.14	36,36,37,38	0
4	DMS	C	8001	4/4	0.94	0.11	34,38,38,40	0
4	DMS	B	8002	4/4	0.94	0.16	42,44,47,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	DMS	C	8009	4/4	0.94	0.13	49,50,51,51	0
2	MG	D	3001	1/1	0.95	0.13	47,47,47,47	0
3	NA	C	3102	1/1	0.95	0.06	23,23,23,23	0
4	DMS	C	8004	4/4	0.95	0.12	27,27,27,28	0
4	DMS	B	8010	4/4	0.95	0.12	47,48,49,49	0
4	DMS	D	8002	4/4	0.95	0.16	40,43,45,47	0
4	DMS	D	8004	4/4	0.95	0.10	24,25,25,28	0
3	NA	A	3103	1/1	0.95	0.10	35,35,35,35	0
4	DMS	D	8006	4/4	0.95	0.11	30,32,33,35	0
2	MG	D	3002	1/1	0.95	0.12	31,31,31,31	0
4	DMS	D	8008	4/4	0.95	0.12	57,57,58,58	0
4	DMS	D	8009	4/4	0.95	0.20	62,62,63,65	0
4	DMS	C	8008	4/4	0.95	0.16	29,29,30,30	0
4	DMS	D	8001	4/4	0.96	0.09	37,38,38,39	0
3	NA	B	3102	1/1	0.96	0.05	18,18,18,18	0
4	DMS	D	8003	4/4	0.96	0.12	26,30,31,38	0
4	DMS	A	8004	4/4	0.96	0.10	41,44,44,44	0
3	NA	A	3102	1/1	0.96	0.07	15,15,15,15	0
3	NA	C	3101	1/1	0.97	0.05	30,30,30,30	0
4	DMS	A	8002	4/4	0.98	0.10	27,29,29,35	0
3	NA	B	3101	1/1	0.98	0.06	24,24,24,24	0
3	NA	D	3102	1/1	0.98	0.07	15,15,15,15	0
4	DMS	A	8001	4/4	0.98	0.07	31,34,34,37	0

6.5 Other polymers

There are no such residues in this entry.