



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 08:08 AM UTC

PDB ID : 3VD9 / pdb_00003vd9
Title : E. coli (lacZ) beta-galactosidase (N460S) in complex with IPTG
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 : 2.05 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

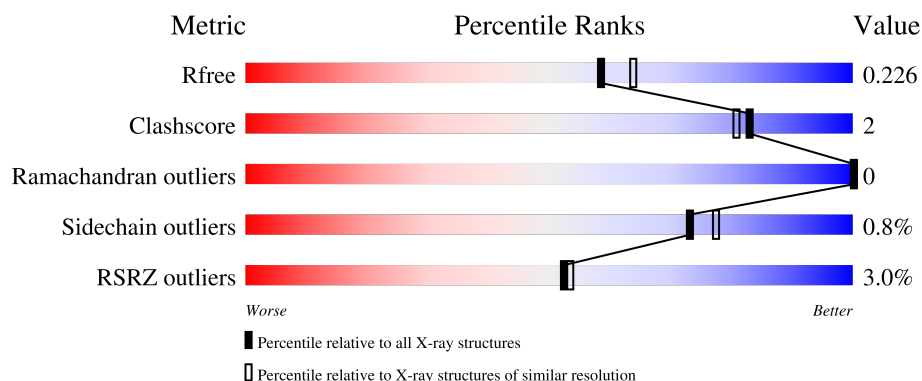
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

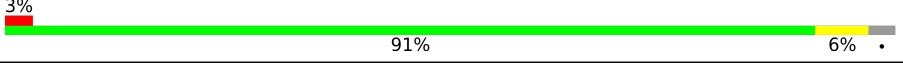
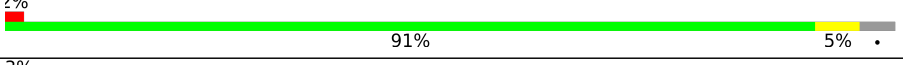
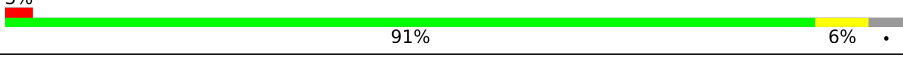
The reported resolution of this entry is 2.05 Å.

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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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 6 unique types of molecules in this entry. The entry contains 36969 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
			8155	5158	1444	1515	38			
1	B	1016	Total	C	N	O	S	0	0	0
			8162	5163	1445	1516	38			
1	C	1015	Total	C	N	O	S	0	0	0
			8155	5158	1444	1515	38			
1	D	1015	Total	C	N	O	S	0	0	0
			8155	5158	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	SER	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	SER	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	SER	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	SER	ASN	engineered mutation	UNP P00722

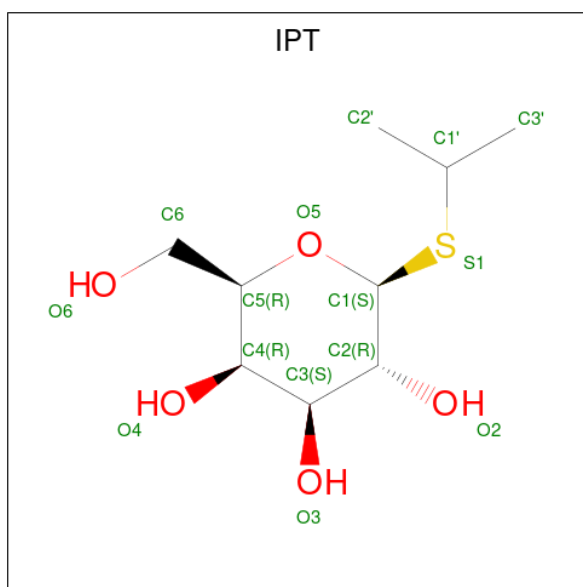
- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		
2	B	3	Total	Mg	0	0
			3	3		
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).

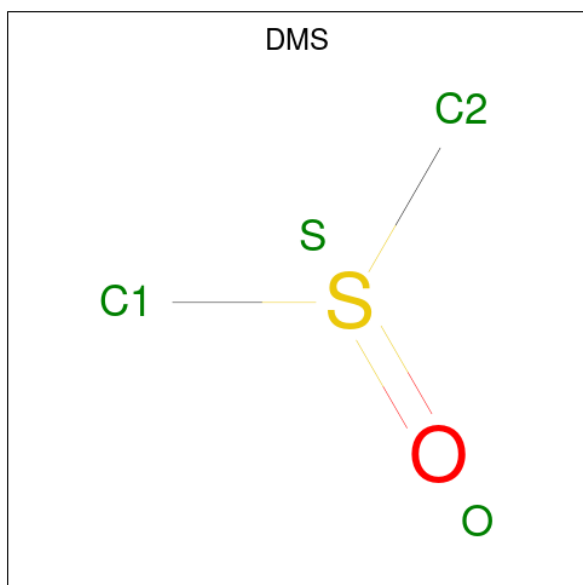
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	Na	0	0
			4	4		
3	B	5	Total	Na	0	0
			5	5		
3	C	4	Total	Na	0	0
			4	4		
3	D	3	Total	Na	0	0
			3	3		

- Molecule 4 is 1-methylethyl 1-thio-beta-D-galactopyranoside (CCD ID: IPT) (formula: C₉H₁₈O₅S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			15	9	5	1		
4	B	1	Total	C	O	S	0	0
			15	9	5	1		
4	C	1	Total	C	O	S	0	0
			15	9	5	1		
4	D	1	Total	C	O	S	0	0
			15	9	5	1		

- Molecule 5 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



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

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

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

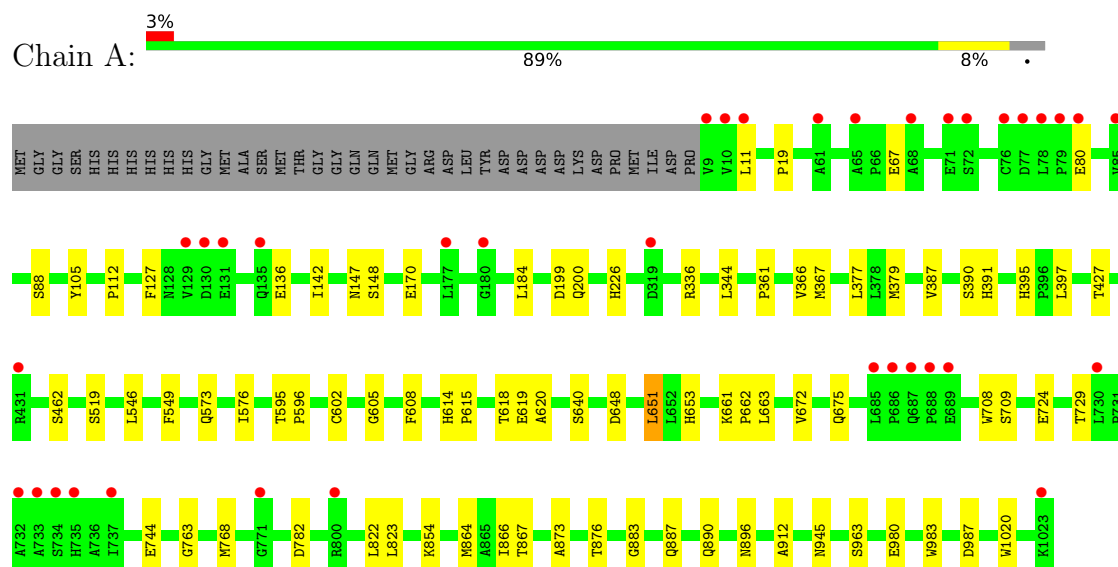
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	918	Total 918	O 918	0	0
6	B	1063	Total 1063	O 1063	0	0
6	C	1099	Total 1099	O 1099	0	0
6	D	921	Total 921	O 921	0	0

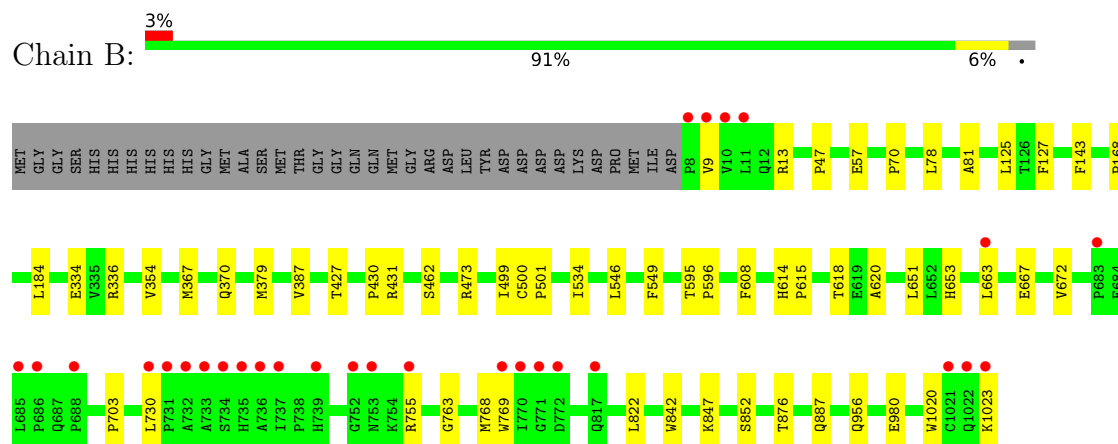
3 Residue-property plots [i](#)

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

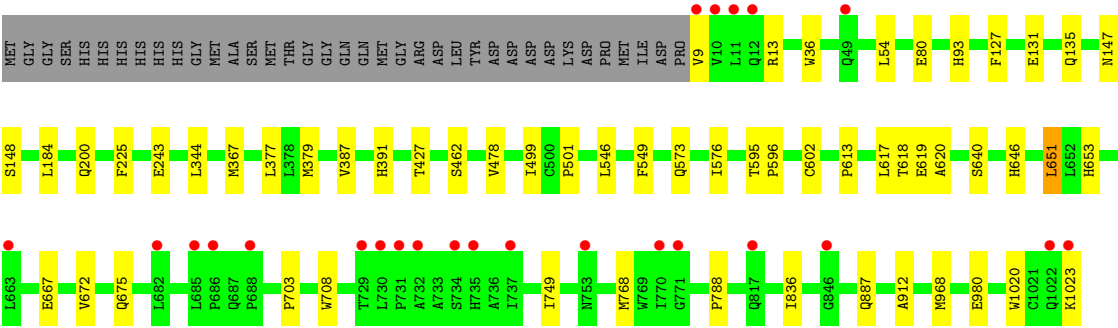


• Molecule 1: Beta-galactosidase

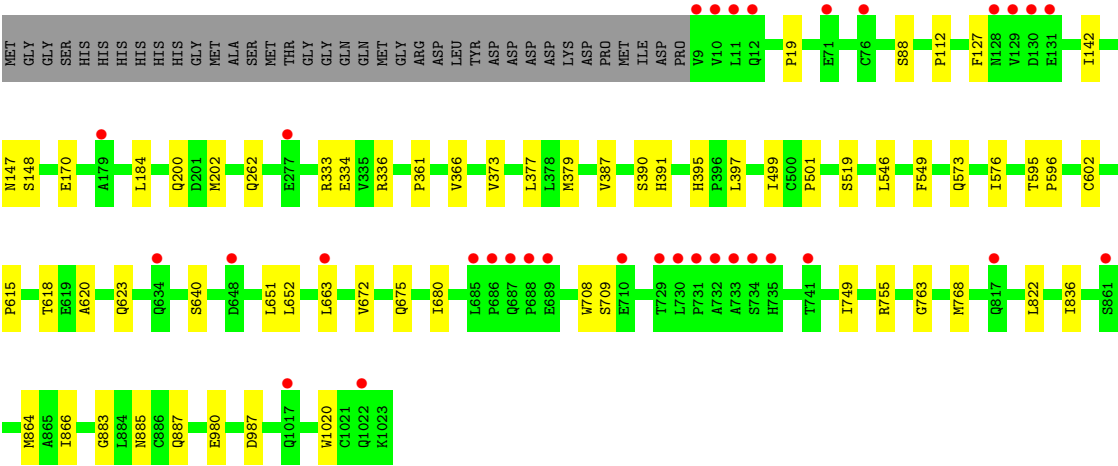
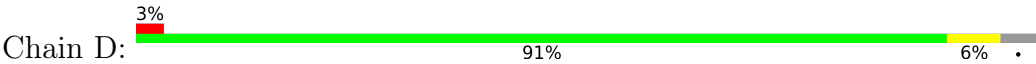


• Molecule 1: Beta-galactosidase





• 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	151.54Å 162.55Å 203.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	86.23 – 2.05 86.23 – 2.05	Depositor EDS
% Data completeness (in resolution range)	100.0 (86.23-2.05) 88.1 (86.23-2.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.05Å)	Xtrriage
Refinement program	REFMAC 5.5.0109, CNS	Depositor
R, R_{free}	0.169 , 0.223 0.174 , 0.226	Depositor DCC
R_{free} test set	3916 reflections (1.42%)	wwPDB-VP
Wilson B-factor (Å ²)	17.0	Xtrriage
Anisotropy	0.073	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	36969	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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 39.66 % 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 3.0702e-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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.47	0/8397	0.70	0/11455
1	B	0.49	0/8405	0.70	0/11466
1	C	0.49	0/8397	0.70	0/11455
1	D	0.47	0/8397	0.70	0/11455
All	All	0.48	0/33596	0.70	0/45831

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	8155	0	7752	48	0
1	B	8162	0	7760	35	0
1	C	8155	0	7752	35	0
1	D	8155	0	7752	34	0
2	A	2	0	0	0	0
2	B	3	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	5	0	0	0	0
3	C	4	0	0	0	0
3	D	3	0	0	0	0
4	A	15	0	17	0	0
4	B	15	0	17	0	0
4	C	15	0	17	0	0
4	D	15	0	17	0	0
5	A	52	0	78	2	0
5	B	80	0	120	7	0
5	C	68	0	102	7	0
5	D	56	0	84	5	0
6	A	918	0	0	4	0
6	B	1063	0	0	2	0
6	C	1099	0	0	5	0
6	D	921	0	0	0	0
All	All	36969	0	31468	152	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 (152) 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:395:HIS:HD2	1:D:397:LEU:H	1.24	0.81
1:C:80:GLU:HG3	6:C:4543:HOH:O	1.84	0.77
1:C:618:THR:HG21	6:C:4337:HOH:O	1.87	0.74
1:A:142:ILE:HG12	1:A:170:GLU:HG2	1.73	0.70
5:C:8008:DMS:H11	6:C:4793:HOH:O	1.94	0.68
1:D:887:GLN:NE2	1:D:980:GLU:O	2.25	0.67
1:D:395:HIS:CD2	1:D:397:LEU:H	2.11	0.67
5:C:8004:DMS:H21	6:C:4423:HOH:O	1.95	0.66
1:C:651:LEU:HD23	1:C:703:PRO:HG3	1.79	0.65
1:B:379:MET:HE1	1:B:387:VAL:HB	1.79	0.64
1:B:768:MET:HE1	1:B:1020:TRP:CZ2	2.34	0.63
1:C:887:GLN:NE2	1:C:980:GLU:O	2.31	0.63
1:A:127:PHE:HE1	1:A:184:LEU:HG	1.64	0.63
1:A:768:MET:HE1	1:A:1020:TRP:CZ2	2.34	0.63
1:B:127:PHE:CE1	1:B:184:LEU:HG	2.35	0.62
1:A:651:LEU:HD11	1:A:653:HIS:CE1	2.37	0.60
1:D:379:MET:HE1	1:D:387:VAL:HB	1.84	0.60
1:A:127:PHE:CE1	1:A:184:LEU:HG	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:ILE:HG12	1:D:170:GLU:HG2	1.84	0.59
1:B:127:PHE:HE1	1:B:184:LEU:HG	1.66	0.59
1:C:377:LEU:HD22	1:C:708:TRP:HA	1.83	0.59
1:A:887:GLN:NE2	1:A:980:GLU:O	2.35	0.58
1:C:54:LEU:HA	5:C:8008:DMS:H12	1.86	0.58
1:B:431:ARG:HG3	6:B:4813:HOH:O	2.03	0.57
1:D:127:PHE:CE1	1:D:184:LEU:HG	2.38	0.57
5:B:8004:DMS:O	5:B:8013:DMS:H21	2.05	0.57
1:A:226:HIS:HD2	6:A:4847:HOH:O	1.88	0.56
1:C:127:PHE:CE1	1:C:184:LEU:HG	2.42	0.55
1:A:708:TRP:CZ2	5:A:8003:DMS:H23	2.41	0.55
1:B:653:HIS:CE1	1:B:667:GLU:HG2	2.41	0.54
1:A:768:MET:HE1	1:A:1020:TRP:CH2	2.43	0.54
1:B:47:PRO:HD3	5:B:8018:DMS:H12	1.91	0.53
1:A:648:ASP:HB3	6:A:4612:HOH:O	2.07	0.53
1:A:377:LEU:HD22	1:A:708:TRP:HA	1.90	0.53
1:B:615:PRO:O	1:B:618:THR:HG22	2.09	0.53
1:D:127:PHE:HE1	1:D:184:LEU:HG	1.74	0.53
1:C:200:GLN:HG2	1:C:391:HIS:HB2	1.90	0.53
1:D:623:GLN:HA	5:D:8002:DMS:O	2.09	0.52
1:D:499:ILE:HG22	1:D:501:PRO:HD3	1.91	0.52
1:D:19:PRO:HD3	1:D:112:PRO:HB3	1.91	0.51
1:C:576:ILE:HD11	6:C:4243:HOH:O	2.10	0.50
1:A:395:HIS:HD2	1:A:397:LEU:H	1.59	0.50
1:D:373:VAL:O	1:D:377:LEU:HG	2.12	0.50
1:B:887:GLN:NE2	1:B:980:GLU:O	2.44	0.50
1:A:651:LEU:HD12	1:A:651:LEU:C	2.37	0.50
1:B:57:GLU:O	5:B:8014:DMS:H12	2.12	0.49
1:D:361:PRO:HB2	1:D:576:ILE:HG12	1.93	0.49
1:C:549:PHE:CE2	1:C:620:ALA:HA	2.47	0.49
1:D:615:PRO:O	1:D:618:THR:HG22	2.12	0.49
1:D:708:TRP:CZ2	5:D:8003:DMS:H23	2.48	0.49
1:B:499:ILE:HG22	1:B:501:PRO:HD3	1.96	0.48
1:A:763:GLY:HA3	1:A:822:LEU:HD13	1.95	0.48
1:C:768:MET:HE1	1:C:1020:TRP:CZ2	2.49	0.48
1:A:708:TRP:CE3	1:A:709:SER:HB3	2.48	0.48
1:C:499:ILE:HG22	1:C:501:PRO:HD3	1.95	0.48
1:C:768:MET:HE1	1:C:1020:TRP:CH2	2.49	0.48
1:D:763:GLY:HA3	1:D:822:LEU:HD13	1.96	0.48
1:A:19:PRO:HD3	1:A:112:PRO:HB3	1.95	0.48
1:A:88:SER:HA	1:A:366:VAL:HG21	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:GLU:HB2	6:A:4656:HOH:O	2.14	0.47
1:C:36:TRP:HH2	5:C:8004:DMS:H12	1.79	0.47
1:D:147:ASN:HA	1:D:148:SER:HA	1.70	0.47
1:A:147:ASN:HA	1:A:148:SER:HA	1.66	0.47
1:D:768:MET:HE1	1:D:1020:TRP:CZ2	2.49	0.47
1:A:864:MET:HE3	1:A:866:ILE:HD11	1.97	0.47
1:C:640:SER:O	1:C:675:GLN:HA	2.15	0.47
1:C:708:TRP:CZ2	5:C:8003:DMS:H23	2.49	0.47
1:D:708:TRP:CE3	1:D:709:SER:HB3	2.50	0.47
1:A:595:THR:HA	1:A:596:PRO:C	2.41	0.46
1:A:549:PHE:CE2	1:A:620:ALA:HA	2.51	0.46
1:D:573:GLN:HB2	1:D:602:CYS:O	2.16	0.45
1:A:573:GLN:HB2	1:A:602:CYS:O	2.16	0.45
1:A:11:LEU:HD11	1:A:67:GLU:HG3	1.98	0.45
1:A:619:GLU:HA	1:A:912:ALA:HB2	1.99	0.45
1:A:873:ALA:O	1:A:876:THR:HG22	2.17	0.45
1:A:615:PRO:O	1:A:618:THR:HG22	2.17	0.45
1:B:549:PHE:CE2	1:B:620:ALA:HA	2.51	0.45
1:A:640:SER:O	1:A:675:GLN:HA	2.17	0.45
1:A:823:LEU:HB3	1:B:730:LEU:HD11	1.99	0.45
1:C:131:GLU:O	1:C:135:GLN:HG3	2.16	0.45
1:B:125:LEU:HD21	5:B:8008:DMS:H13	1.99	0.45
1:C:613:PRO:HB3	1:C:617:LEU:HD23	1.99	0.45
1:D:749:ILE:HD11	1:D:836:ILE:HD11	1.98	0.45
1:A:379:MET:HE1	1:A:387:VAL:HB	1.98	0.44
1:B:595:THR:HA	1:B:596:PRO:C	2.42	0.44
1:D:377:LEU:HD22	1:D:708:TRP:HA	1.98	0.44
1:D:334:GLU:OE1	1:D:336:ARG:NH1	2.46	0.44
1:D:390:SER:HA	1:D:391:HIS:HA	1.80	0.44
1:A:427:THR:HG21	1:A:462:SER:HB3	1.98	0.44
1:C:147:ASN:HA	1:C:148:SER:HA	1.72	0.44
1:A:782:ASP:OD1	1:A:854:LYS:NZ	2.44	0.44
1:A:896:ASN:HB3	1:A:945:ASN:HB2	1.99	0.44
1:D:640:SER:O	1:D:675:GLN:HA	2.18	0.44
1:A:336:ARG:HB3	5:A:8007:DMS:H23	2.00	0.44
1:B:354:VAL:HG12	1:B:379:MET:HE3	2.00	0.43
1:B:9:VAL:HG22	1:C:9:VAL:HG22	1.99	0.43
1:B:370:GLN:HB2	6:B:4798:HOH:O	2.17	0.43
1:B:427:THR:HG21	1:B:462:SER:HB3	2.00	0.43
1:B:755:ARG:HB3	1:B:769:TRP:HB2	1.99	0.43
1:B:334:GLU:OE1	1:B:336:ARG:NH1	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:842:TRP:HZ3	1:B:852:SER:HB3	1.82	0.43
1:D:88:SER:HA	1:D:366:VAL:HG21	2.00	0.43
1:A:200:GLN:HG2	1:A:391:HIS:HB2	1.99	0.43
1:B:143:PHE:O	1:B:168:PRO:HA	2.18	0.43
1:B:430:PRO:HD3	5:B:8012:DMS:H23	2.00	0.43
1:B:367:MET:HA	1:B:367:MET:HE2	1.99	0.43
1:D:652:LEU:HD22	1:D:680:ILE:HD13	2.01	0.43
1:A:724:GLU:O	1:B:847:LYS:NZ	2.52	0.43
1:B:763:GLY:HA3	1:B:822:LEU:HD13	2.01	0.43
1:C:225:PHE:HA	1:C:243:GLU:O	2.19	0.43
1:A:605:GLY:O	1:A:614:HIS:ND1	2.52	0.42
1:A:661:LYS:HA	1:A:662:PRO:HD3	1.92	0.42
1:D:596:PRO:HG2	5:D:8014:DMS:H11	2.01	0.42
1:B:81:ALA:HA	5:B:8016:DMS:H12	2.00	0.42
1:B:651:LEU:HD23	1:B:703:PRO:HG3	2.01	0.42
1:C:595:THR:HA	1:C:596:PRO:C	2.45	0.42
1:D:595:THR:HA	1:D:596:PRO:C	2.44	0.42
1:C:749:ILE:HD11	1:C:836:ILE:HD11	2.01	0.42
1:C:788:PRO:HD2	1:C:968:MET:HB2	2.01	0.42
1:C:576:ILE:HG12	5:C:8011:DMS:C2	2.50	0.42
1:C:367:MET:HE2	1:C:367:MET:HA	2.02	0.42
1:C:573:GLN:HB2	1:C:602:CYS:O	2.20	0.42
1:D:883:GLY:HA3	1:D:987:ASP:HA	2.02	0.42
1:D:333:ARG:NH1	5:D:8001:DMS:H12	2.34	0.42
1:A:651:LEU:HD11	1:A:653:HIS:ND1	2.35	0.41
1:B:13:ARG:HG3	1:C:13:ARG:CZ	2.50	0.41
1:D:19:PRO:HD3	1:D:112:PRO:CB	2.49	0.41
5:D:8004:DMS:O	5:D:8013:DMS:H11	2.20	0.41
1:C:619:GLU:HA	1:C:912:ALA:HB2	2.02	0.41
5:B:8012:DMS:H13	1:C:478:VAL:HG22	2.02	0.41
1:A:367:MET:HA	1:A:367:MET:HE2	2.03	0.41
1:C:427:THR:HG21	1:C:462:SER:HB3	2.03	0.41
1:D:864:MET:HE3	1:D:866:ILE:HD11	2.02	0.41
1:A:390:SER:HA	1:A:391:HIS:HA	1.81	0.41
1:B:70:PRO:HG2	1:B:78:LEU:HD21	2.02	0.41
1:B:651:LEU:HD11	1:B:653:HIS:NE2	2.36	0.41
1:C:93:HIS:CG	5:C:8013:DMS:H12	2.55	0.41
1:A:854:LYS:HA	1:A:867:THR:O	2.20	0.41
1:A:883:GLY:HA3	1:A:987:ASP:HA	2.03	0.41
1:B:500:CYS:HA	1:B:534:ILE:O	2.21	0.41
1:C:653:HIS:CE1	1:C:667:GLU:HG2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:TYR:CE1	1:A:199:ASP:HB2	2.56	0.40
1:B:473:ARG:HA	1:B:473:ARG:HD2	1.93	0.40
1:C:127:PHE:HE1	1:C:184:LEU:HG	1.85	0.40
1:A:361:PRO:HB2	1:A:576:ILE:HG12	2.03	0.40
1:A:890:GLN:HB2	6:A:4728:HOH:O	2.21	0.40
1:D:200:GLN:HB2	1:D:202:MET:HE2	2.02	0.40
1:B:608:PHE:CE2	1:B:614:HIS:HD2	2.40	0.40
1:C:379:MET:HE1	1:C:387:VAL:HB	2.04	0.40
1:D:549:PHE:CE2	1:D:620:ALA:HA	2.57	0.40
1:A:608:PHE:CE2	1:A:614:HIS:HD2	2.39	0.40
1:A:963:SER:HB3	1:A:983:TRP:CE2	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	1013/1052 (96%)	978 (96%)	35 (4%)	0	100	100
1	B	1014/1052 (96%)	982 (97%)	32 (3%)	0	100	100
1	C	1013/1052 (96%)	980 (97%)	33 (3%)	0	100	100
1	D	1013/1052 (96%)	982 (97%)	31 (3%)	0	100	100
All	All	4053/4208 (96%)	3922 (97%)	131 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	868/898 (97%)	859 (99%)	9 (1%)	68	72
1	B	869/898 (97%)	863 (99%)	6 (1%)	76	80
1	C	868/898 (97%)	862 (99%)	6 (1%)	76	80
1	D	868/898 (97%)	860 (99%)	8 (1%)	70	75
All	All	3473/3592 (97%)	3444 (99%)	29 (1%)	73	77

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	GLU
1	A	344	LEU
1	A	519	SER
1	A	546	LEU
1	A	651	LEU
1	A	663	LEU
1	A	672	VAL
1	A	729	THR
1	A	744	GLU
1	B	546	LEU
1	B	663	LEU
1	B	672	VAL
1	B	876	THR
1	B	956	GLN
1	B	1023	LYS
1	C	344	LEU
1	C	546	LEU
1	C	646	HIS
1	C	651	LEU
1	C	672	VAL
1	C	1023	LYS
1	D	262	GLN
1	D	519	SER
1	D	546	LEU
1	D	651	LEU
1	D	663	LEU
1	D	672	VAL
1	D	755	ARG
1	D	885	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	50	GLN
1	A	221	GLN
1	A	445	GLN
1	A	485	GLN
1	A	624	GLN
1	A	702	GLN
1	A	843	GLN
1	B	23	GLN
1	B	50	GLN
1	B	266	GLN
1	B	485	GLN
1	B	702	GLN
1	B	804	ASN
1	B	824	GLN
1	B	903	GLN
1	B	950	GLN
1	C	363	HIS
1	C	365	GLN
1	C	485	GLN
1	C	634	GLN
1	C	702	GLN
1	C	885	ASN
1	C	950	GLN
1	D	50	GLN
1	D	128	ASN
1	D	395	HIS
1	D	445	GLN
1	D	634	GLN
1	D	702	GLN
1	D	843	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 93 ligands modelled in this entry, 25 are monoatomic - leaving 68 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
5	DMS	B	8018	-	3,3,3	2.70	1 (33%)	3,3,3	0.56	0
5	DMS	C	8004	-	3,3,3	2.68	1 (33%)	3,3,3	0.58	0
5	DMS	B	8015	-	3,3,3	2.67	1 (33%)	3,3,3	0.50	0
5	DMS	D	8009	-	3,3,3	2.75	1 (33%)	3,3,3	0.49	0
5	DMS	A	8005	-	3,3,3	2.63	1 (33%)	3,3,3	0.41	0
5	DMS	B	8012	-	3,3,3	2.68	1 (33%)	3,3,3	0.60	0
4	IPT	B	2001	3	15,15,15	0.71	0	18,21,21	0.85	0
5	DMS	D	8011	-	3,3,3	2.69	1 (33%)	3,3,3	0.50	0
5	DMS	C	8005	-	3,3,3	1.91	1 (33%)	3,3,3	0.29	0
5	DMS	A	8008	-	3,3,3	2.66	1 (33%)	3,3,3	0.52	0
5	DMS	A	8006	-	3,3,3	2.71	1 (33%)	3,3,3	0.67	0
5	DMS	B	8019	-	3,3,3	2.74	1 (33%)	3,3,3	0.38	0
5	DMS	C	8006	-	3,3,3	2.67	1 (33%)	3,3,3	0.53	0
5	DMS	D	8008	-	3,3,3	2.09	1 (33%)	3,3,3	0.45	0
5	DMS	C	8003	-	3,3,3	1.98	1 (33%)	3,3,3	0.41	0
5	DMS	C	8017	-	3,3,3	2.67	1 (33%)	3,3,3	0.54	0
5	DMS	C	8009	-	3,3,3	2.66	1 (33%)	3,3,3	0.58	0
4	IPT	D	2001	3	15,15,15	0.68	0	18,21,21	0.81	0
5	DMS	C	8008	-	3,3,3	2.70	1 (33%)	3,3,3	0.68	0
5	DMS	C	8010	-	3,3,3	2.76	1 (33%)	3,3,3	0.60	0
5	DMS	C	8002	-	3,3,3	2.66	1 (33%)	3,3,3	0.61	0
5	DMS	B	8008	-	3,3,3	2.72	1 (33%)	3,3,3	0.56	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DMS	B	8017	-	3,3,3	2.73	1 (33%)	3,3,3	0.57	0
5	DMS	A	8007	-	3,3,3	2.73	1 (33%)	3,3,3	0.65	0
4	IPT	A	2001	3	15,15,15	0.66	0	18,21,21	0.71	0
5	DMS	A	8011	-	3,3,3	2.71	1 (33%)	3,3,3	0.55	0
5	DMS	B	8010	-	3,3,3	2.63	1 (33%)	3,3,3	0.40	0
5	DMS	C	8014	-	3,3,3	2.71	1 (33%)	3,3,3	0.56	0
5	DMS	C	8015	-	3,3,3	2.70	1 (33%)	3,3,3	0.60	0
5	DMS	D	8006	-	3,3,3	2.70	1 (33%)	3,3,3	0.52	0
5	DMS	C	8011	-	3,3,3	2.62	1 (33%)	3,3,3	0.36	0
5	DMS	C	8012	-	3,3,3	2.70	1 (33%)	3,3,3	0.66	0
4	IPT	C	2001	3	15,15,15	0.65	0	18,21,21	0.81	0
5	DMS	A	8009	-	3,3,3	2.68	1 (33%)	3,3,3	0.45	0
5	DMS	B	8007	-	3,3,3	2.71	1 (33%)	3,3,3	0.52	0
5	DMS	B	8002	-	3,3,3	2.60	1 (33%)	3,3,3	0.50	0
5	DMS	A	8013	-	3,3,3	2.71	1 (33%)	3,3,3	0.55	0
5	DMS	D	8001	-	3,3,3	2.67	1 (33%)	3,3,3	0.84	0
5	DMS	B	8005	-	3,3,3	2.65	1 (33%)	3,3,3	0.38	0
5	DMS	D	8002	-	3,3,3	2.58	1 (33%)	3,3,3	0.32	0
5	DMS	D	8004	-	3,3,3	2.65	1 (33%)	3,3,3	0.46	0
5	DMS	D	8005	-	3,3,3	2.63	1 (33%)	3,3,3	0.31	0
5	DMS	C	8007	-	3,3,3	2.70	1 (33%)	3,3,3	0.51	0
5	DMS	C	8001	-	3,3,3	2.65	1 (33%)	3,3,3	0.62	0
5	DMS	C	8013	-	3,3,3	2.67	1 (33%)	3,3,3	0.66	0
5	DMS	A	8001	-	3,3,3	1.98	1 (33%)	3,3,3	0.60	0
5	DMS	B	8016	-	3,3,3	2.63	1 (33%)	3,3,3	0.58	0
5	DMS	B	8009	-	3,3,3	2.75	1 (33%)	3,3,3	0.57	0
5	DMS	D	8007	-	3,3,3	2.71	1 (33%)	3,3,3	0.68	0
5	DMS	B	8001	-	3,3,3	2.62	1 (33%)	3,3,3	0.52	0
5	DMS	B	8004	-	3,3,3	2.62	1 (33%)	3,3,3	0.63	0
5	DMS	A	8003	-	3,3,3	2.02	1 (33%)	3,3,3	0.50	0
5	DMS	B	8006	-	3,3,3	2.71	1 (33%)	3,3,3	0.69	0
5	DMS	D	8014	-	3,3,3	2.71	1 (33%)	3,3,3	0.53	0
5	DMS	B	8003	-	3,3,3	1.93	1 (33%)	3,3,3	0.53	0
5	DMS	A	8010	-	3,3,3	2.70	1 (33%)	3,3,3	0.56	0
5	DMS	B	8013	-	3,3,3	2.64	1 (33%)	3,3,3	0.49	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DMS	B	8020	-	3,3,3	2.70	1 (33%)	3,3,3	0.33	0
5	DMS	D	8013	-	3,3,3	2.72	1 (33%)	3,3,3	0.60	0
5	DMS	A	8002	-	3,3,3	2.64	1 (33%)	3,3,3	0.51	0
5	DMS	D	8003	-	3,3,3	2.03	1 (33%)	3,3,3	0.38	0
5	DMS	A	8004	-	3,3,3	2.67	1 (33%)	3,3,3	0.54	0
5	DMS	D	8010	-	3,3,3	2.62	1 (33%)	3,3,3	0.70	0
5	DMS	B	8011	-	3,3,3	1.99	1 (33%)	3,3,3	0.42	0
5	DMS	A	8012	-	3,3,3	2.72	1 (33%)	3,3,3	0.79	0
5	DMS	C	8016	-	3,3,3	2.77	1 (33%)	3,3,3	0.60	0
5	DMS	B	8014	-	3,3,3	2.70	1 (33%)	3,3,3	0.72	0
5	DMS	D	8012	-	3,3,3	2.73	1 (33%)	3,3,3	0.56	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
4	IPT	D	2001	3	-	3/6/26/26	0/1/1/1
4	IPT	C	2001	3	-	4/6/26/26	0/1/1/1
4	IPT	B	2001	3	-	3/6/26/26	0/1/1/1
4	IPT	A	2001	3	-	3/6/26/26	0/1/1/1

All (64) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	8016	DMS	O-S	4.63	1.80	1.50
5	C	8010	DMS	O-S	4.63	1.80	1.50
5	B	8009	DMS	O-S	4.63	1.80	1.50
5	D	8009	DMS	O-S	4.60	1.80	1.50
5	B	8019	DMS	O-S	4.60	1.80	1.50
5	A	8007	DMS	O-S	4.59	1.80	1.50
5	B	8017	DMS	O-S	4.58	1.80	1.50
5	D	8013	DMS	O-S	4.58	1.80	1.50
5	B	8020	DMS	O-S	4.57	1.80	1.50
5	A	8013	DMS	O-S	4.57	1.80	1.50
5	B	8006	DMS	O-S	4.57	1.80	1.50
5	B	8008	DMS	O-S	4.56	1.80	1.50
5	A	8011	DMS	O-S	4.56	1.80	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	8012	DMS	O-S	4.56	1.80	1.50
5	A	8012	DMS	O-S	4.55	1.80	1.50
5	D	8014	DMS	O-S	4.55	1.80	1.50
5	C	8012	DMS	O-S	4.55	1.80	1.50
5	C	8014	DMS	O-S	4.55	1.80	1.50
5	B	8018	DMS	O-S	4.54	1.80	1.50
5	C	8008	DMS	O-S	4.54	1.80	1.50
5	A	8010	DMS	O-S	4.54	1.80	1.50
5	A	8006	DMS	O-S	4.54	1.80	1.50
5	D	8007	DMS	O-S	4.54	1.80	1.50
5	C	8015	DMS	O-S	4.54	1.80	1.50
5	B	8007	DMS	O-S	4.54	1.80	1.50
5	B	8014	DMS	O-S	4.53	1.80	1.50
5	C	8007	DMS	O-S	4.52	1.80	1.50
5	D	8006	DMS	O-S	4.51	1.80	1.50
5	B	8012	DMS	O-S	4.51	1.80	1.50
5	C	8004	DMS	O-S	4.51	1.80	1.50
5	D	8011	DMS	O-S	4.50	1.79	1.50
5	A	8004	DMS	O-S	4.50	1.79	1.50
5	A	8009	DMS	O-S	4.50	1.79	1.50
5	C	8006	DMS	O-S	4.50	1.79	1.50
5	B	8015	DMS	O-S	4.49	1.79	1.50
5	C	8013	DMS	O-S	4.48	1.79	1.50
5	C	8002	DMS	O-S	4.47	1.79	1.50
5	A	8008	DMS	O-S	4.47	1.79	1.50
5	C	8017	DMS	O-S	4.46	1.79	1.50
5	D	8001	DMS	O-S	4.46	1.79	1.50
5	B	8013	DMS	O-S	4.46	1.79	1.50
5	C	8009	DMS	O-S	4.46	1.79	1.50
5	D	8004	DMS	O-S	4.45	1.79	1.50
5	C	8001	DMS	O-S	4.45	1.79	1.50
5	A	8002	DMS	O-S	4.44	1.79	1.50
5	B	8005	DMS	O-S	4.44	1.79	1.50
5	B	8010	DMS	O-S	4.43	1.79	1.50
5	B	8016	DMS	O-S	4.43	1.79	1.50
5	D	8005	DMS	O-S	4.42	1.79	1.50
5	C	8011	DMS	O-S	4.41	1.79	1.50
5	B	8004	DMS	O-S	4.40	1.79	1.50
5	A	8005	DMS	O-S	4.40	1.79	1.50
5	B	8001	DMS	O-S	4.38	1.79	1.50
5	D	8010	DMS	O-S	4.38	1.79	1.50
5	B	8002	DMS	O-S	4.37	1.79	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	8002	DMS	O-S	4.34	1.78	1.50
5	D	8008	DMS	O-S	3.56	1.73	1.50
5	D	8003	DMS	O-S	3.44	1.72	1.50
5	A	8003	DMS	O-S	3.42	1.72	1.50
5	B	8011	DMS	O-S	3.39	1.72	1.50
5	C	8003	DMS	O-S	3.37	1.72	1.50
5	A	8001	DMS	O-S	3.36	1.72	1.50
5	B	8003	DMS	O-S	3.28	1.71	1.50
5	C	8005	DMS	O-S	3.24	1.71	1.50

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2001	IPT	C2'-C1'-S1-C1
4	A	2001	IPT	C3'-C1'-S1-C1
4	B	2001	IPT	C2'-C1'-S1-C1
4	B	2001	IPT	C3'-C1'-S1-C1
4	C	2001	IPT	C2'-C1'-S1-C1
4	C	2001	IPT	C3'-C1'-S1-C1
4	D	2001	IPT	C2'-C1'-S1-C1
4	D	2001	IPT	C3'-C1'-S1-C1
4	C	2001	IPT	O5-C5-C6-O6
4	D	2001	IPT	O5-C5-C6-O6
4	A	2001	IPT	O5-C5-C6-O6
4	B	2001	IPT	O5-C5-C6-O6
4	C	2001	IPT	C4-C5-C6-O6

There are no ring outliers.

20 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	8018	DMS	1	0
5	C	8004	DMS	2	0
5	B	8012	DMS	2	0
5	C	8003	DMS	1	0
5	C	8008	DMS	2	0
5	B	8008	DMS	1	0
5	A	8007	DMS	1	0
5	C	8011	DMS	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	8001	DMS	1	0
5	D	8002	DMS	1	0
5	D	8004	DMS	1	0
5	C	8013	DMS	1	0
5	B	8016	DMS	1	0
5	B	8004	DMS	1	0
5	A	8003	DMS	1	0
5	D	8014	DMS	1	0
5	B	8013	DMS	1	0
5	D	8013	DMS	1	0
5	D	8003	DMS	1	0
5	B	8014	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%)	0.01	36 (3%)	47	48	14, 26, 48, 78	0
1	B	1016/1052 (96%)	-0.26	29 (2%)	53	55	12, 21, 46, 80	0
1	C	1015/1052 (96%)	-0.33	24 (2%)	59	61	11, 21, 43, 75	0
1	D	1015/1052 (96%)	-0.02	33 (3%)	49	49	15, 26, 47, 82	0
All	All	4061/4208 (96%)	-0.15	122 (3%)	52	53	11, 24, 47, 82	0

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	770	ILE	5.1
1	B	11	LEU	5.0
1	C	10	VAL	4.9
1	A	10	VAL	4.8
1	A	730	LEU	4.6
1	B	731	PRO	4.6
1	D	730	LEU	4.4
1	B	733	ALA	4.4
1	B	9	VAL	4.3
1	C	11	LEU	4.3
1	D	10	VAL	4.3
1	A	9	VAL	4.2
1	A	11	LEU	4.1
1	B	730	LEU	4.1
1	B	771	GLY	4.1
1	A	733	ALA	4.1
1	C	9	VAL	4.0
1	C	686	PRO	3.9
1	B	1022	GLN	3.9
1	C	1022	GLN	3.9
1	C	1023	LYS	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	685	LEU	3.8
1	B	10	VAL	3.6
1	B	663	LEU	3.6
1	B	685	LEU	3.6
1	D	735	HIS	3.5
1	B	8	PRO	3.5
1	D	733	ALA	3.5
1	A	735	HIS	3.5
1	C	730	LEU	3.4
1	A	686	PRO	3.4
1	C	731	PRO	3.4
1	D	11	LEU	3.4
1	B	732	ALA	3.4
1	C	770	ILE	3.4
1	B	772	ASP	3.3
1	D	129	VAL	3.3
1	A	688	PRO	3.3
1	D	9	VAL	3.3
1	A	732	ALA	3.2
1	B	688	PRO	3.2
1	D	732	ALA	3.2
1	D	76	CYS	3.2
1	C	735	HIS	3.0
1	A	76	CYS	3.0
1	A	79	PRO	3.0
1	B	734	SER	3.0
1	D	741	THR	3.0
1	B	686	PRO	3.0
1	B	1023	LYS	3.0
1	D	685	LEU	2.9
1	B	752	GLY	2.9
1	D	131	GLU	2.9
1	D	686	PRO	2.9
1	A	771	GLY	2.9
1	D	689	GLU	2.9
1	D	734	SER	2.8
1	D	688	PRO	2.8
1	A	689	GLU	2.7
1	C	685	LEU	2.7
1	D	12	GLN	2.7
1	D	731	PRO	2.7
1	D	710	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	735	HIS	2.6
1	C	688	PRO	2.6
1	C	753	ASN	2.6
1	D	130	ASP	2.6
1	A	78	LEU	2.6
1	A	135	GLN	2.5
1	D	687	GLN	2.5
1	A	131	GLU	2.5
1	C	771	GLY	2.5
1	B	753	ASN	2.5
1	A	68	ALA	2.5
1	A	687	GLN	2.5
1	C	12	GLN	2.5
1	B	736	ALA	2.5
1	D	179	ALA	2.5
1	B	737	ILE	2.5
1	D	729	THR	2.4
1	C	663	LEU	2.4
1	A	129	VAL	2.4
1	B	755	ARG	2.4
1	C	729	THR	2.4
1	A	77	ASP	2.4
1	C	734	SER	2.4
1	B	769	TRP	2.4
1	A	65	ALA	2.3
1	A	130	ASP	2.3
1	A	734	SER	2.3
1	D	128	ASN	2.3
1	A	72	SER	2.3
1	D	861	SER	2.3
1	D	648	ASP	2.3
1	C	846	GLY	2.3
1	A	80	GLU	2.2
1	C	49	GLN	2.2
1	C	737	ILE	2.2
1	B	739	HIS	2.2
1	C	817	GLN	2.2
1	A	85	VAL	2.2
1	A	180	GLY	2.2
1	D	634	GLN	2.2
1	D	1017	GLN	2.2
1	A	800	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	737	ILE	2.1
1	D	71	GLU	2.1
1	A	1023	LYS	2.1
1	C	682	LEU	2.1
1	D	663	LEU	2.1
1	D	277	GLU	2.1
1	A	319	ASP	2.1
1	D	1022	GLN	2.1
1	A	71	GLU	2.1
1	B	817	GLN	2.1
1	A	431	ARG	2.1
1	C	732	ALA	2.1
1	B	1021	CYS	2.0
1	D	817	GLN	2.0
1	B	683	PRO	2.0
1	A	61	ALA	2.0
1	A	177	LEU	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
5	DMS	D	8006	4/4	0.81	0.22	50,53,53,58	0
5	DMS	D	8009	4/4	0.81	0.24	71,72,73,74	0
5	DMS	B	8007	4/4	0.82	0.23	47,49,52,53	0
5	DMS	A	8006	4/4	0.83	0.29	63,63,63,67	0
5	DMS	A	8012	4/4	0.84	0.22	51,54,56,59	0
5	DMS	B	8012	4/4	0.84	0.25	54,56,58,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	DMS	D	8007	4/4	0.86	0.26	61,61,62,62	0
5	DMS	C	8007	4/4	0.86	0.21	51,52,52,55	0
5	DMS	D	8013	4/4	0.86	0.22	54,60,60,60	0
5	DMS	A	8010	4/4	0.87	0.27	61,64,64,65	0
5	DMS	A	8011	4/4	0.87	0.18	58,60,61,63	0
3	NA	C	3104	1/1	0.87	0.12	35,35,35,35	0
5	DMS	C	8010	4/4	0.87	0.22	50,52,53,54	0
5	DMS	D	8014	4/4	0.87	0.21	67,68,68,68	0
5	DMS	B	8014	4/4	0.88	0.17	42,48,48,53	0
5	DMS	A	8013	4/4	0.88	0.21	66,66,68,69	0
5	DMS	A	8007	4/4	0.88	0.20	57,58,58,59	0
3	NA	B	3104	1/1	0.88	0.11	39,39,39,39	0
5	DMS	D	8001	4/4	0.89	0.19	38,42,47,49	0
5	DMS	B	8017	4/4	0.89	0.17	58,58,59,62	0
5	DMS	C	8013	4/4	0.89	0.21	56,57,57,58	0
5	DMS	D	8002	4/4	0.90	0.22	38,41,42,48	0
5	DMS	B	8016	4/4	0.90	0.18	47,47,49,55	0
5	DMS	B	8006	4/4	0.90	0.20	46,50,52,52	0
5	DMS	C	8014	4/4	0.90	0.19	63,64,65,68	0
5	DMS	D	8010	4/4	0.90	0.19	48,52,53,53	0
5	DMS	C	8015	4/4	0.90	0.20	54,55,55,59	0
3	NA	B	3105	1/1	0.90	0.22	27,27,27,27	1
5	DMS	C	8004	4/4	0.91	0.17	48,48,50,56	0
3	NA	A	3105	1/1	0.91	0.10	30,30,30,30	1
5	DMS	D	8012	4/4	0.91	0.15	38,44,47,53	0
5	DMS	B	8020	4/4	0.91	0.20	54,55,56,56	0
5	DMS	C	8016	4/4	0.91	0.15	56,56,57,58	0
5	DMS	D	8004	4/4	0.92	0.17	42,45,49,50	0
5	DMS	A	8004	4/4	0.92	0.17	45,47,51,51	0
5	DMS	C	8009	4/4	0.92	0.18	54,55,57,57	0
5	DMS	B	8004	4/4	0.92	0.15	31,34,41,42	0
5	DMS	B	8018	4/4	0.93	0.14	47,48,49,50	0
5	DMS	C	8006	4/4	0.93	0.16	49,50,50,52	0
5	DMS	B	8005	4/4	0.93	0.16	40,41,43,44	0
5	DMS	C	8008	4/4	0.93	0.14	42,48,49,51	0
3	NA	B	3103	1/1	0.94	0.08	29,29,29,29	0
3	NA	D	3101	1/1	0.94	0.07	26,26,26,26	0
5	DMS	B	8019	4/4	0.94	0.13	43,44,46,47	0
3	NA	A	3103	1/1	0.94	0.07	39,39,39,39	0
5	DMS	B	8015	4/4	0.94	0.14	44,45,47,49	0
2	MG	D	3002	1/1	0.94	0.07	25,25,25,25	0
5	DMS	C	8002	4/4	0.95	0.10	20,31,33,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NA	D	3102	1/1	0.95	0.05	22,22,22,22	0
5	DMS	A	8009	4/4	0.95	0.14	41,47,47,48	0
4	IPT	D	2001	15/15	0.95	0.08	26,28,33,34	0
5	DMS	A	8003	4/4	0.95	0.11	40,43,43,45	0
3	NA	B	3102	1/1	0.95	0.06	22,22,22,22	0
2	MG	A	3002	1/1	0.95	0.07	27,27,27,27	0
4	IPT	C	2001	15/15	0.96	0.08	20,26,32,35	0
3	NA	C	3102	1/1	0.96	0.04	20,20,20,20	0
5	DMS	A	8002	4/4	0.96	0.12	33,36,36,38	0
2	MG	B	3007	1/1	0.96	0.05	47,47,47,47	0
2	MG	D	3001	1/1	0.96	0.04	23,23,23,23	0
5	DMS	D	8003	4/4	0.96	0.12	42,43,43,44	0
2	MG	B	3002	1/1	0.96	0.08	19,19,19,19	0
3	NA	D	3103	1/1	0.96	0.07	32,32,32,32	0
5	DMS	B	8008	4/4	0.96	0.12	44,46,46,49	0
5	DMS	B	8009	4/4	0.96	0.10	39,39,41,41	0
5	DMS	B	8011	4/4	0.96	0.11	46,47,47,49	0
5	DMS	D	8011	4/4	0.96	0.09	31,36,41,42	0
5	DMS	A	8008	4/4	0.96	0.11	42,45,46,47	0
4	IPT	A	2001	15/15	0.96	0.08	17,23,29,32	0
4	IPT	B	2001	15/15	0.96	0.08	16,25,30,32	0
5	DMS	B	8002	4/4	0.97	0.08	22,30,30,31	0
5	DMS	B	8003	4/4	0.97	0.11	35,38,38,39	0
5	DMS	D	8005	4/4	0.97	0.10	34,34,37,39	0
5	DMS	C	8011	4/4	0.97	0.09	26,28,28,29	0
5	DMS	C	8003	4/4	0.97	0.09	38,40,40,40	0
5	DMS	D	8008	4/4	0.97	0.11	51,52,52,53	0
2	MG	C	3002	1/1	0.97	0.06	18,18,18,18	0
5	DMS	C	8005	4/4	0.97	0.10	31,33,34,36	0
3	NA	C	3103	1/1	0.97	0.06	33,33,33,33	0
5	DMS	C	8017	4/4	0.97	0.09	28,35,36,36	0
3	NA	A	3102	1/1	0.97	0.04	18,18,18,18	0
5	DMS	B	8013	4/4	0.97	0.08	30,32,35,37	0
5	DMS	C	8001	4/4	0.98	0.07	22,23,28,29	0
5	DMS	A	8001	4/4	0.98	0.09	30,36,37,37	0
5	DMS	B	8010	4/4	0.98	0.09	32,38,38,38	0
5	DMS	A	8005	4/4	0.98	0.08	30,32,34,35	0
5	DMS	C	8012	4/4	0.98	0.10	37,39,39,43	0
5	DMS	B	8001	4/4	0.98	0.08	23,26,27,30	0
3	NA	A	3101	1/1	0.98	0.04	23,23,23,23	0
2	MG	C	3001	1/1	0.98	0.02	17,17,17,17	0
2	MG	A	3001	1/1	0.99	0.02	21,21,21,21	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NA	C	3101	1/1	0.99	0.02	15,15,15,15	0
2	MG	B	3001	1/1	0.99	0.02	16,16,16,16	0
3	NA	B	3101	1/1	0.99	0.04	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.