



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 01:23 PM UTC

PDB ID : 3T2P / pdb_00003t2p
Title : E. coli (lacZ) beta-galactosidase (S796D) in complex with IPTG
Authors : Jancewicz, L.J.; Wheatley, R.W.; Sutendra, G.; Lee, M.; Fraser, M.; Huber, R.E.
Deposited on : 2011-07-22
Resolution : 2.60 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

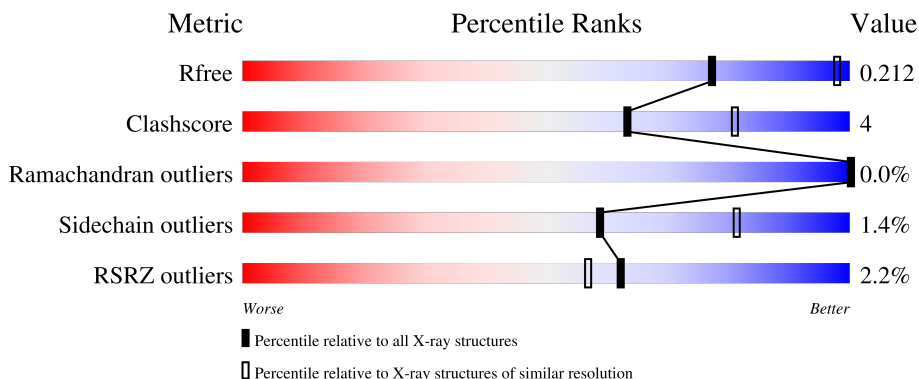
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	 2% 85% 11% .
1	B	1052	 2% 86% 11% .
1	C	1052	 2% 85% 12% .
1	D	1052	 2% 86% 11% .

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 34272 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	1014	Total	C	N	O	S	0	0	0
			8152	5155	1444	1515	38			
1	B	1015	Total	C	N	O	S	0	0	0
			8159	5160	1445	1516	38			
1	C	1015	Total	C	N	O	S	0	0	0
			8159	5160	1445	1516	38			
1	D	1015	Total	C	N	O	S	0	0	0
			8159	5160	1445	1516	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	796	ASP	SER	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	796	ASP	SER	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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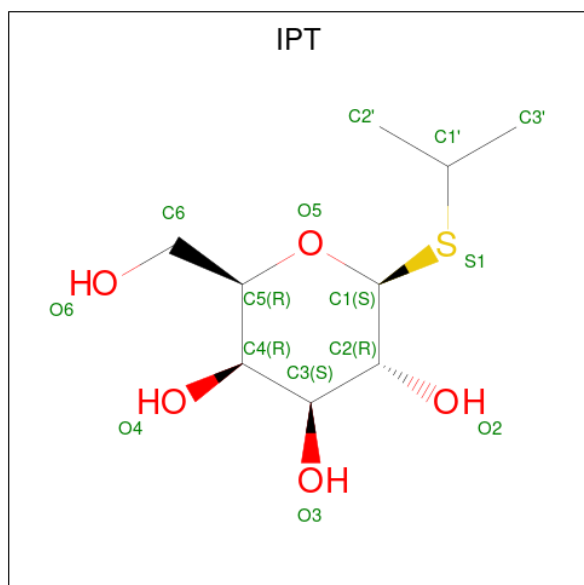
Chain	Residue	Modelled	Actual	Comment	Reference
C	1	ASP	-	expression tag	UNP P00722
C	2	LYS	-	expression tag	UNP P00722
C	3	ASP	-	expression tag	UNP P00722
C	4	PRO	-	expression tag	UNP P00722
C	5	MET	-	expression tag	UNP P00722
C	6	ILE	-	expression tag	UNP P00722
C	7	ASP	-	expression tag	UNP P00722
C	8	PRO	-	expression tag	UNP P00722
C	796	ASP	SER	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	796	ASP	SER	engineered mutation	UNP P00722

- Molecule 2 is 1-methylethyl 1-thio-beta-D-galactopyranoside (CCD ID: IPT) (formula: $C_9H_{18}O_5S$).



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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

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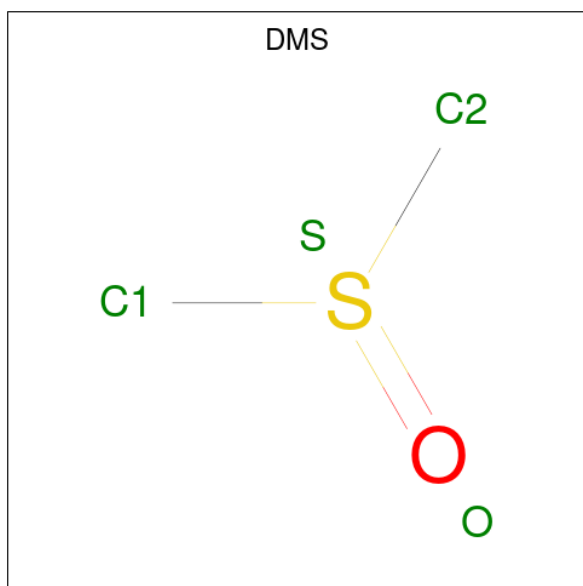
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	2	Total	Mg	0	0
			2	2		
3	D	2	Total	Mg	0	0
			2	2		

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

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

- Molecule 5 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



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

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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	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

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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	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
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	C	O	S	0	0
			4	2	1	1		
5	C	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		
5	D	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	348	Total	O	0	0
			348	348		

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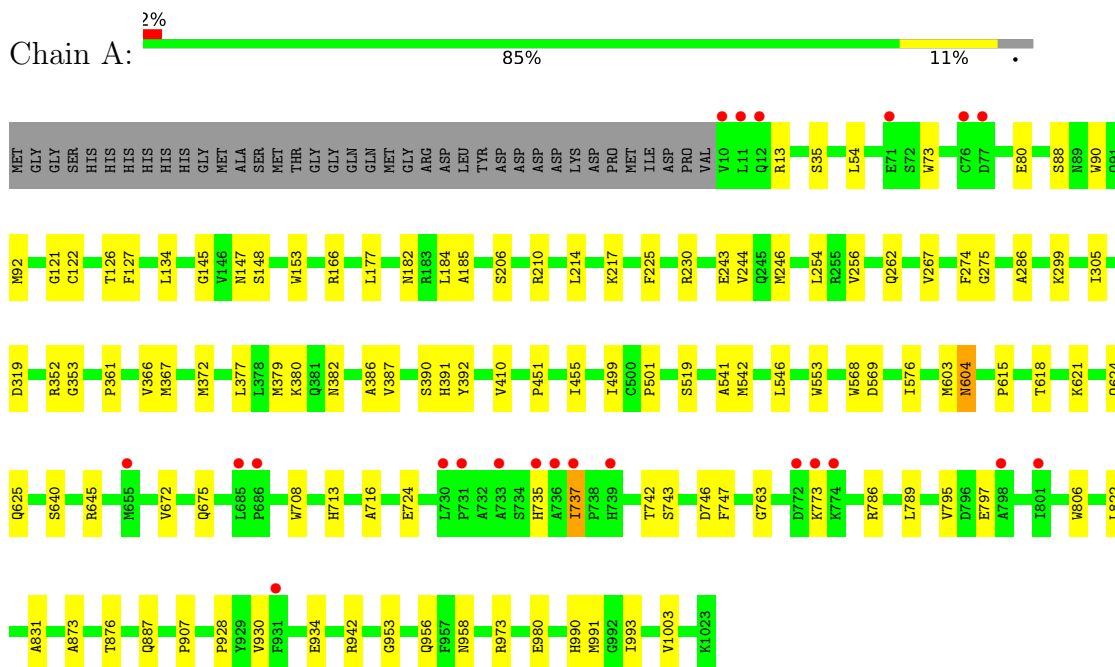
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	270	Total 270	O 270	0	0
6	C	322	Total 322	O 322	0	0
6	D	373	Total 373	O 373	0	0

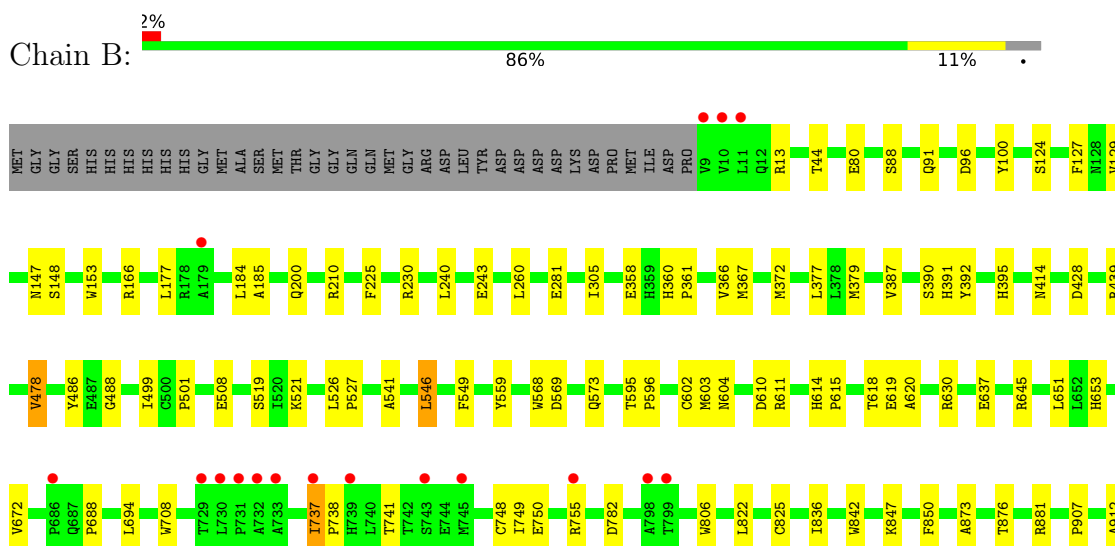
3 Residue-property plots

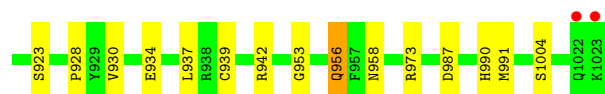
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Beta-galactosidase

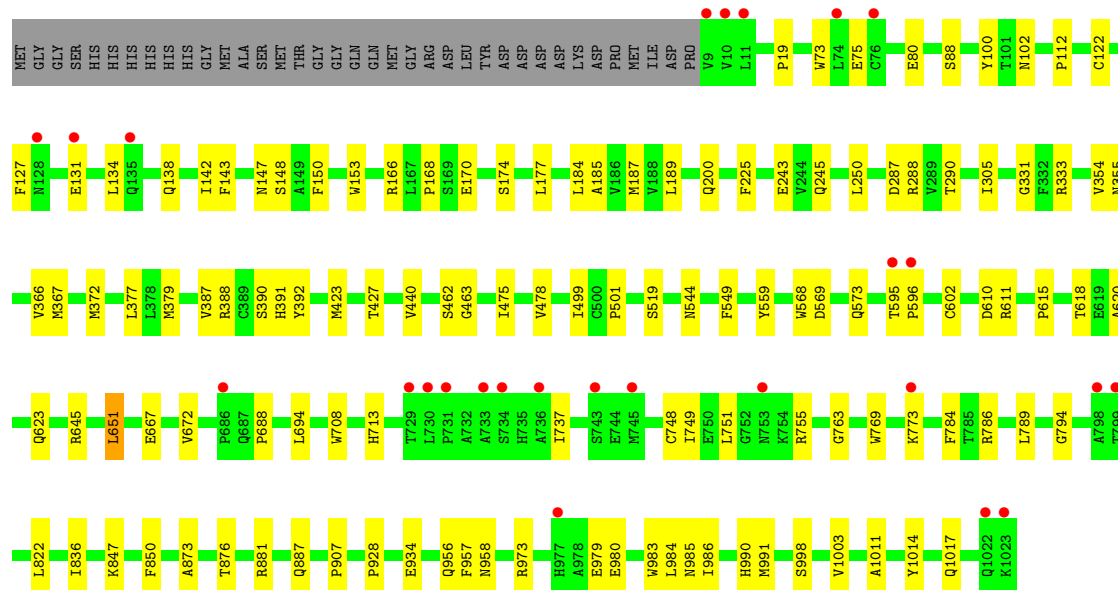
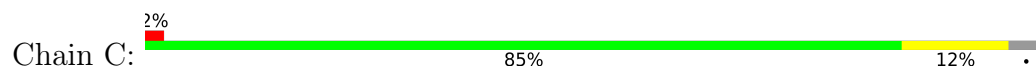


• Molecule 1: Beta-galactosidase

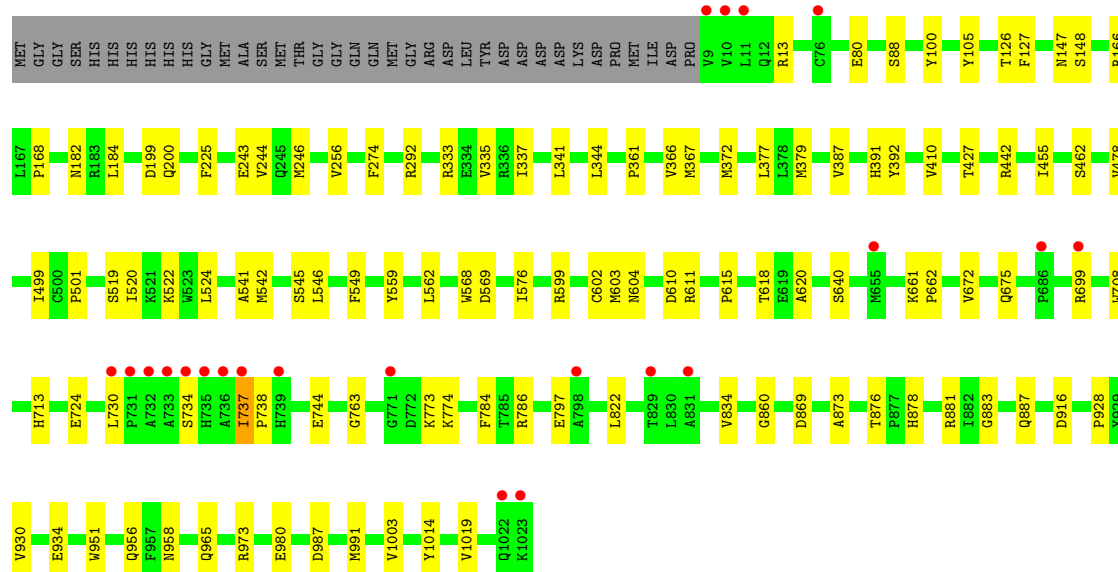
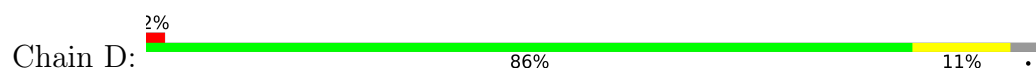




• Molecule 1: Beta-galactosidase



• 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	150.87Å 166.17Å 200.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	74.67 – 2.60 74.67 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (74.67-2.60) 98.7 (74.67-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.01 (at 2.62Å)	Xtrriage
Refinement program	REFMAC 5.5.0109, CNS	Depositor
R, R_{free}	0.184 , 0.242 (Not available) , 0.212	Depositor DCC
R_{free} test set	2160 reflections (1.41%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtrriage
Anisotropy	0.050	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.1	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	34272	wwPDB-VP
Average B, all atoms (Å ²)	52.0	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 43.59 % 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 1.7241e-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: DMS, NA, IPT, MG

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	0.44	0/8394	0.72	0/11451
1	B	0.44	0/8401	0.71	0/11461
1	C	0.43	0/8401	0.71	0/11461
1	D	0.44	0/8401	0.72	0/11461
All	All	0.44	0/33597	0.71	0/45834

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	8152	0	7743	68	0
1	B	8159	0	7752	61	0
1	C	8159	0	7752	70	0
1	D	8159	0	7752	61	0
2	A	15	0	17	0	0
2	B	15	0	17	0	0
2	C	15	0	18	0	0
2	D	15	0	17	0	0
3	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	3	0	0	0	0
4	B	2	0	0	0	0
4	C	3	0	0	0	0
4	D	2	0	0	0	0
5	A	56	0	84	5	0
5	B	60	0	90	3	0
5	C	72	0	108	6	0
5	D	64	0	96	5	0
6	A	348	0	0	3	0
6	B	270	0	0	4	0
6	C	322	0	0	3	0
6	D	373	0	0	4	0
All	All	34272	0	31446	258	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 4.

All (258) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:MET:HE1	1:A:254:LEU:HD13	1.31	1.08
1:B:367:MET:HB3	1:B:372:MET:HE2	1.70	0.73
1:D:377:LEU:HD22	1:D:708:TRP:HA	1.73	0.70
1:C:367:MET:HB3	1:C:372:MET:HE2	1.73	0.70
1:A:451:PRO:HA	5:A:8001:DMS:H23	1.74	0.69
1:D:786:ARG:HB2	1:D:934:GLU:HG3	1.76	0.67
1:C:245:GLN:HG2	1:C:288:ARG:HG2	1.76	0.66
1:A:786:ARG:HB2	1:A:934:GLU:CG	2.27	0.65
1:C:379:MET:HE1	1:C:387:VAL:HB	1.79	0.64
1:D:88:SER:HA	1:D:366:VAL:HG21	1.80	0.64
1:D:367:MET:HB3	1:D:372:MET:HE2	1.79	0.63
1:D:379:MET:HE1	1:D:387:VAL:HB	1.79	0.63
1:C:847:LYS:NZ	1:D:724:GLU:O	2.32	0.63
1:A:379:MET:HE1	1:A:387:VAL:HB	1.81	0.62
1:A:246:MET:HE3	1:A:274:PHE:CE2	2.35	0.62
1:A:377:LEU:HD22	1:A:708:TRP:HA	1.83	0.60
1:A:786:ARG:HB2	1:A:934:GLU:HG3	1.81	0.60
1:A:615:PRO:O	1:A:618:THR:HG22	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:GLY:HA2	1:A:286:ALA:HA	1.84	0.59
1:D:878:HIS:HD2	6:D:4286:HOH:O	1.85	0.59
1:C:713:HIS:HD2	6:C:4287:HOH:O	1.83	0.59
1:C:153:TRP:HB2	1:C:185:ALA:HB3	1.84	0.59
1:A:367:MET:HB3	1:A:372:MET:HE2	1.85	0.58
1:C:290:THR:O	5:C:8012:DMS:H22	2.04	0.58
1:A:305:ILE:HD11	1:A:645:ARG:HB3	1.85	0.58
1:D:873:ALA:O	1:D:876:THR:HG22	2.03	0.58
1:D:615:PRO:O	1:D:618:THR:HG22	2.04	0.57
1:A:724:GLU:O	1:B:847:LYS:NZ	2.38	0.57
1:C:934:GLU:OE1	1:C:958:ASN:ND2	2.38	0.57
1:C:377:LEU:HD22	1:C:708:TRP:HA	1.88	0.56
1:A:887:GLN:NE2	1:A:980:GLU:O	2.37	0.56
1:D:699:ARG:NH2	5:D:8013:DMS:H12	2.21	0.56
1:B:88:SER:HA	1:B:366:VAL:HG21	1.87	0.56
1:D:887:GLN:NE2	1:D:980:GLU:O	2.35	0.55
1:A:640:SER:O	1:A:675:GLN:HA	2.06	0.55
1:B:614:HIS:HB3	6:B:4068:HOH:O	2.06	0.54
1:C:763:GLY:HA3	1:C:822:LEU:HD13	1.90	0.54
1:C:615:PRO:O	1:C:618:THR:HG22	2.07	0.53
1:D:610:ASP:O	1:D:611:ARG:HB2	2.09	0.53
1:A:451:PRO:HA	5:A:8001:DMS:C2	2.38	0.53
1:D:499:ILE:HG22	1:D:501:PRO:HD3	1.90	0.53
1:D:603:MET:HE1	1:D:930:VAL:HB	1.91	0.53
1:A:763:GLY:HA3	1:A:822:LEU:HD13	1.91	0.53
1:C:499:ILE:HG22	1:C:501:PRO:HD3	1.90	0.53
1:B:508:GLU:OE2	5:B:8007:DMS:H11	2.09	0.52
1:B:541:ALA:HB3	1:B:604:ASN:O	2.10	0.52
1:C:305:ILE:HD11	1:C:645:ARG:HB3	1.91	0.52
1:A:147:ASN:HB3	1:A:206:SER:HA	1.92	0.52
1:C:573:GLN:HB2	1:C:602:CYS:O	2.10	0.52
1:A:873:ALA:O	1:A:876:THR:HG22	2.10	0.51
1:C:623:GLN:HA	5:C:8002:DMS:H12	1.92	0.51
1:C:651:LEU:HD11	1:C:667:GLU:HB3	1.91	0.51
1:A:499:ILE:HG22	1:A:501:PRO:HD3	1.92	0.51
1:B:478:VAL:HG22	5:C:8015:DMS:H13	1.92	0.51
1:D:934:GLU:OE1	1:D:958:ASN:ND2	2.43	0.51
1:B:615:PRO:O	1:B:618:THR:HG22	2.11	0.51
1:B:428:ASP:O	5:B:8012:DMS:H12	2.10	0.51
1:C:127:PHE:HE1	1:C:184:LEU:HG	1.76	0.51
1:A:166:ARG:HG3	1:A:392:TYR:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:542:MET:HA	1:A:604:ASN:HA	1.93	0.51
1:D:127:PHE:CE1	1:D:184:LEU:HG	2.46	0.50
5:D:8002:DMS:H22	6:D:4034:HOH:O	2.10	0.50
1:D:427:THR:HG21	1:D:462:SER:HB3	1.93	0.50
1:C:755:ARG:HB3	1:C:769:TRP:HB2	1.93	0.50
1:B:942:ARG:HA	1:B:953:GLY:O	2.12	0.50
1:C:142:ILE:HG12	1:C:170:GLU:HG2	1.94	0.49
1:B:782:ASP:HB2	1:B:842:TRP:CZ2	2.47	0.49
1:A:134:LEU:HD21	1:A:177:LEU:HB2	1.94	0.49
1:B:546:LEU:HA	6:B:4056:HOH:O	2.12	0.49
1:D:730:LEU:HD12	1:D:730:LEU:H	1.78	0.49
1:A:907:PRO:HG2	1:A:990:HIS:O	2.12	0.49
1:C:134:LEU:HD21	1:C:177:LEU:HB2	1.94	0.49
1:C:568:TRP:CD2	1:C:569:ASP:HB3	2.48	0.49
1:C:928:PRO:HB2	1:C:973:ARG:HH11	1.78	0.49
1:C:786:ARG:HB2	1:C:934:GLU:HG3	1.93	0.49
1:A:361:PRO:HB2	1:A:576:ILE:HG12	1.95	0.48
1:B:603:MET:HE1	1:B:930:VAL:HB	1.95	0.48
1:C:127:PHE:CE1	1:C:184:LEU:HG	2.48	0.48
1:D:361:PRO:HB2	1:D:576:ILE:HG12	1.95	0.48
1:A:934:GLU:OE1	1:A:958:ASN:ND2	2.46	0.48
1:C:427:THR:HG21	1:C:462:SER:HB3	1.96	0.48
1:B:750:GLU:HG3	1:B:755:ARG:HG3	1.95	0.48
1:B:939:CYS:HA	1:B:956:GLN:HB3	1.96	0.48
1:C:595:THR:HA	1:C:596:PRO:C	2.39	0.48
1:B:100:TYR:CE1	1:B:602:CYS:HB3	2.48	0.47
1:B:305:ILE:HD11	1:B:645:ARG:HB3	1.94	0.47
1:D:100:TYR:CE1	1:D:602:CYS:HB3	2.49	0.47
1:B:822:LEU:HD11	1:B:825:CYS:HB2	1.95	0.47
1:A:737:ILE:HD12	1:A:831:ALA:O	2.15	0.47
1:B:934:GLU:OE1	1:B:958:ASN:ND2	2.47	0.47
1:C:887:GLN:NE2	1:C:980:GLU:O	2.45	0.47
1:C:907:PRO:HG2	1:C:990:HIS:O	2.15	0.47
1:A:127:PHE:CE1	1:A:184:LEU:HG	2.50	0.47
1:B:881:ARG:HE	1:B:987:ASP:CG	2.23	0.47
1:C:991:MET:HE3	1:C:1003:VAL:HG21	1.96	0.47
1:C:200:GLN:HG2	1:C:391:HIS:HB2	1.97	0.47
1:D:127:PHE:HE1	1:D:184:LEU:HG	1.80	0.47
1:A:88:SER:HA	1:A:366:VAL:HG21	1.98	0.46
1:C:688:PRO:HG3	1:C:694:LEU:HD21	1.97	0.46
5:B:8012:DMS:H13	1:C:478:VAL:HG22	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:200:GLN:HG2	1:D:391:HIS:HB2	1.97	0.46
1:A:942:ARG:HA	1:A:953:GLY:O	2.16	0.46
1:B:200:GLN:HG2	1:B:391:HIS:HB2	1.97	0.46
1:B:390:SER:HA	1:B:391:HIS:HA	1.76	0.46
1:A:367:MET:HB3	1:A:372:MET:CE	2.45	0.46
1:A:991:MET:HE2	1:A:1003:VAL:HG21	1.98	0.46
1:A:806:TRP:CE2	1:A:991:MET:HE1	2.51	0.46
1:C:786:ARG:NH2	1:C:991:MET:HE2	2.31	0.46
1:A:127:PHE:HE1	1:A:184:LEU:HG	1.81	0.46
1:B:595:THR:HA	1:B:596:PRO:C	2.40	0.46
1:C:225:PHE:HA	1:C:243:GLU:O	2.16	0.46
1:C:549:PHE:CE2	1:C:620:ALA:HA	2.51	0.45
1:D:734:SER:OG	1:D:860:GLY:HA3	2.16	0.45
1:B:521:LYS:HG2	1:B:559:TYR:CZ	2.51	0.45
1:D:786:ARG:NH2	1:D:991:MET:HE2	2.32	0.45
1:B:907:PRO:HG2	1:B:990:HIS:O	2.17	0.45
1:C:143:PHE:O	1:C:168:PRO:HA	2.17	0.45
1:C:354:VAL:CG1	1:C:379:MET:HE3	2.46	0.45
1:D:549:PHE:CE2	1:D:620:ALA:HA	2.51	0.45
1:C:984:LEU:HD21	1:C:986:ILE:HD11	1.99	0.45
1:A:380:LYS:O	5:A:8003:DMS:H11	2.17	0.45
1:A:390:SER:HA	1:A:391:HIS:HA	1.83	0.45
1:B:526:LEU:O	1:B:527:PRO:C	2.60	0.45
1:D:166:ARG:HG3	1:D:392:TYR:HB2	1.98	0.45
1:B:873:ALA:O	1:B:876:THR:HG22	2.17	0.45
1:D:225:PHE:HA	1:D:243:GLU:O	2.17	0.45
1:D:524:LEU:HD11	1:D:562:LEU:HG	1.99	0.45
1:B:610:ASP:O	1:B:611:ARG:HB2	2.17	0.45
1:C:786:ARG:HB2	1:C:934:GLU:CG	2.47	0.45
1:D:168:PRO:O	1:D:442:ARG:NH2	2.50	0.45
1:D:568:TRP:CD2	1:D:569:ASP:HB3	2.52	0.45
1:A:928:PRO:HB2	1:A:973:ARG:HH11	1.82	0.44
1:B:499:ILE:HG22	1:B:501:PRO:HD3	1.99	0.44
1:D:246:MET:HG2	1:D:274:PHE:CE1	2.52	0.44
1:D:883:GLY:HA3	1:D:987:ASP:HA	2.00	0.44
1:B:379:MET:HE1	1:B:387:VAL:HB	1.98	0.44
1:A:603:MET:HE1	1:A:930:VAL:HB	1.99	0.44
1:B:166:ARG:HG3	1:B:392:TYR:HB2	1.98	0.44
1:B:549:PHE:CE2	1:B:620:ALA:HA	2.51	0.44
1:C:544:ASN:HB3	1:C:789:LEU:HD22	2.00	0.44
1:C:651:LEU:CD1	1:C:667:GLU:HB3	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:292:ARG:H	5:D:8005:DMS:H11	1.82	0.44
1:B:573:GLN:HB2	1:B:602:CYS:O	2.17	0.44
1:C:333:ARG:NH1	5:C:8001:DMS:H12	2.33	0.44
1:D:699:ARG:HH21	5:D:8013:DMS:H12	1.82	0.44
1:A:625:GLN:HG2	1:A:716:ALA:HA	2.00	0.43
1:D:147:ASN:HA	1:D:148:SER:HA	1.69	0.43
1:D:951:TRP:HA	1:D:1019:VAL:O	2.18	0.43
1:B:127:PHE:CE1	1:B:184:LEU:HG	2.53	0.43
1:B:129:VAL:HG21	1:B:177:LEU:HD22	2.00	0.43
1:B:568:TRP:CD2	1:B:569:ASP:HB3	2.53	0.43
1:D:784:PHE:HA	1:D:881:ARG:O	2.18	0.43
1:A:90:TRP:CZ3	1:A:121:GLY:HA3	2.53	0.43
1:D:559:TYR:HB2	1:D:562:LEU:HD12	1.99	0.43
1:B:651:LEU:HD11	1:B:653:HIS:CE1	2.53	0.43
1:B:737:ILE:HA	1:B:738:PRO:HD3	1.87	0.43
1:C:979:GLU:OE1	1:C:983:TRP:NE1	2.50	0.43
1:B:153:TRP:HB2	1:B:185:ALA:HB3	1.99	0.43
1:D:333:ARG:NH1	5:D:8001:DMS:H12	2.34	0.43
1:D:928:PRO:HB2	1:D:973:ARG:HH11	1.83	0.43
1:A:73:TRP:CE2	1:A:122:CYS:HB3	2.54	0.43
1:C:749:ILE:HD11	1:C:836:ILE:HD11	2.01	0.43
1:D:661:LYS:HA	1:D:662:PRO:HD3	1.90	0.43
1:A:145:GLY:N	1:A:210:ARG:HB2	2.34	0.43
1:A:742:THR:HG23	1:A:747:PHE:HE1	1.84	0.43
1:C:1017:GLN:HB2	6:C:4252:HOH:O	2.18	0.43
1:A:230:ARG:HB3	6:A:4258:HOH:O	2.19	0.43
1:B:749:ILE:HD11	1:B:836:ILE:HD11	2.00	0.43
1:C:873:ALA:O	1:C:876:THR:HG22	2.19	0.43
1:B:13:ARG:HA	1:B:13:ARG:HD3	1.87	0.43
1:B:281:GLU:HB2	1:C:423:MET:CE	2.49	0.43
1:B:414:ASN:O	1:B:439:ARG:HD3	2.19	0.43
1:B:630:ARG:HB2	1:B:637:GLU:HB3	2.01	0.43
1:B:881:ARG:NH2	1:B:934:GLU:OE1	2.42	0.43
1:A:13:ARG:CZ	1:D:13:ARG:HG3	2.49	0.42
1:B:124:SER:HA	1:B:184:LEU:O	2.18	0.42
1:C:100:TYR:CE1	1:C:602:CYS:HB3	2.54	0.42
1:D:335:VAL:HG22	1:D:344:LEU:HD12	2.01	0.42
1:D:991:MET:HE3	1:D:1003:VAL:HG21	2.00	0.42
1:B:741:THR:HB	1:B:748:CYS:HB2	2.01	0.42
5:C:8004:DMS:O	5:C:8017:DMS:H21	2.19	0.42
1:D:520:ILE:HD12	1:D:562:LEU:HD22	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:541:ALA:HB3	1:D:604:ASN:O	2.19	0.42
1:D:640:SER:O	1:D:675:GLN:HA	2.19	0.42
1:D:763:GLY:HA3	1:D:822:LEU:HD13	2.01	0.42
1:C:610:ASP:O	1:C:611:ARG:HB2	2.19	0.42
1:A:126:THR:HA	1:A:182:ASN:O	2.18	0.42
1:A:153:TRP:HB2	1:A:185:ALA:HB3	2.01	0.42
1:A:553:TRP:CZ2	1:A:624:GLN:HG2	2.54	0.42
1:B:619:GLU:HA	1:B:912:ALA:HB2	2.00	0.42
1:B:688:PRO:HG3	1:B:694:LEU:HD21	2.01	0.42
1:D:737:ILE:HA	1:D:738:PRO:HD3	1.89	0.42
1:B:210:ARG:NH2	1:B:358:GLU:OE1	2.47	0.42
1:D:244:VAL:HG21	1:D:256:VAL:HG11	2.01	0.42
1:D:410:VAL:HG22	1:D:455:ILE:HB	2.01	0.42
1:A:713:HIS:HE1	6:A:4130:HOH:O	2.02	0.42
1:B:230:ARG:HD3	6:B:4215:HOH:O	2.18	0.42
1:C:250:LEU:HD21	1:C:287:ASP:HA	2.01	0.42
1:C:440:VAL:HG13	1:C:475:ILE:HD11	2.02	0.42
1:D:126:THR:HA	1:D:182:ASN:O	2.20	0.42
1:D:542:MET:O	1:D:545:SER:HB2	2.18	0.42
1:A:225:PHE:HA	1:A:243:GLU:O	2.20	0.42
6:B:4158:HOH:O	1:C:463:GLY:HA2	2.19	0.42
1:C:784:PHE:HA	1:C:881:ARG:O	2.20	0.42
1:B:372:MET:HE1	1:B:395:HIS:CG	2.54	0.42
1:A:410:VAL:HG22	1:A:455:ILE:HB	2.02	0.42
1:A:806:TRP:CD2	1:A:991:MET:HE1	2.54	0.42
1:D:337:ILE:HA	1:D:341:LEU:O	2.20	0.42
1:D:599:ARG:HH21	1:D:797:GLU:HG2	1.83	0.42
1:A:541:ALA:HB3	1:A:604:ASN:O	2.20	0.42
1:B:806:TRP:CD2	1:B:991:MET:HE1	2.55	0.42
1:C:794:GLY:HA2	1:C:998:SER:O	2.20	0.41
1:A:286:ALA:HB2	5:A:8016:DMS:H22	2.02	0.41
1:C:1011:ALA:HB3	1:C:1014:TYR:CZ	2.55	0.41
1:A:352:ARG:HG2	1:A:553:TRP:CH2	2.55	0.41
1:A:568:TRP:HA	1:A:569:ASP:HA	1.83	0.41
1:B:91:GLN:HG3	1:B:96:ASP:OD1	2.19	0.41
1:A:382:ASN:HA	1:A:621:LYS:HD2	2.01	0.41
1:A:35:SER:HB2	1:A:217:LYS:HD2	2.02	0.41
1:A:743:SER:HB3	1:A:746:ASP:O	2.21	0.41
1:B:486:TYR:CZ	1:B:488:GLY:HA3	2.56	0.41
1:C:19:PRO:HD3	1:C:112:PRO:CB	2.51	0.41
1:B:928:PRO:HB2	1:B:973:ARG:HH11	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:138:GLN:HA	1:C:174:SER:OG	2.20	0.41
1:A:92:MET:HA	1:A:92:MET:HE2	2.02	0.41
1:B:147:ASN:HA	1:B:148:SER:HA	1.67	0.41
1:C:147:ASN:HA	1:C:148:SER:HA	1.71	0.41
1:C:354:VAL:HG12	1:C:379:MET:HE3	2.02	0.41
1:C:166:ARG:HG3	1:C:392:TYR:HB2	2.02	0.41
1:C:331:GLY:O	5:C:8001:DMS:H13	2.21	0.41
1:D:105:TYR:CE1	1:D:199:ASP:HB2	2.56	0.41
1:A:147:ASN:HA	1:A:148:SER:HA	1.61	0.41
1:A:244:VAL:HG21	1:A:256:VAL:HG11	2.03	0.41
1:A:262:GLN:NE2	1:A:299:LYS:HD2	2.35	0.41
1:A:353:GLY:HA2	1:A:386:ALA:O	2.21	0.41
1:A:786:ARG:HB2	1:A:934:GLU:HG2	2.01	0.41
1:B:225:PHE:HA	1:B:243:GLU:O	2.21	0.41
1:B:360:HIS:ND1	1:B:361:PRO:HD2	2.35	0.41
1:C:390:SER:HA	1:C:391:HIS:HA	1.83	0.41
1:A:246:MET:HE3	1:A:274:PHE:CD2	2.56	0.41
1:C:150:PHE:HA	1:C:187:MET:O	2.20	0.41
1:A:13:ARG:HA	1:A:13:ARG:HD3	1.90	0.40
1:B:377:LEU:HD22	1:B:708:TRP:HA	2.02	0.40
1:C:88:SER:HA	1:C:366:VAL:HG21	2.03	0.40
1:C:559:TYR:CE2	1:D:522:LYS:HA	2.56	0.40
1:A:54:LEU:HD11	1:A:214:LEU:HG	2.04	0.40
1:C:355:ASN:OD1	1:C:388:ARG:HD3	2.21	0.40
1:D:713:HIS:HE1	6:D:4166:HOH:O	2.03	0.40
1:A:789:LEU:HD11	1:A:993:ILE:HG22	2.04	0.40
1:C:713:HIS:CD2	6:C:4287:HOH:O	2.67	0.40
1:D:774:LYS:HE3	6:D:4251:HOH:O	2.21	0.40
1:B:240:LEU:HD13	1:B:260:LEU:HD13	2.04	0.40
1:C:73:TRP:CE2	1:C:122:CYS:HB3	2.56	0.40
1:C:166:ARG:HG3	1:C:392:TYR:CB	2.51	0.40
1:C:957:PHE:HA	1:C:985:ASN:O	2.21	0.40
5:A:8003:DMS:H13	6:A:4040:HOH:O	2.21	0.40
1:D:738:PRO:HB2	1:D:834:VAL:HG23	2.04	0.40
1:D:869:ASP:HA	1:D:1014:TYR:O	2.20	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	1012/1052 (96%)	971 (96%)	40 (4%)	1 (0%)	48	70
1	B	1013/1052 (96%)	971 (96%)	42 (4%)	0	100	100
1	C	1013/1052 (96%)	969 (96%)	43 (4%)	1 (0%)	48	70
1	D	1013/1052 (96%)	974 (96%)	39 (4%)	0	100	100
All	All	4051/4208 (96%)	3885 (96%)	164 (4%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	795	VAL
1	C	102	ASN

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	867/898 (96%)	855 (99%)	12 (1%)	59	81
1	B	868/898 (97%)	856 (99%)	12 (1%)	59	81
1	C	868/898 (97%)	855 (98%)	13 (2%)	57	80
1	D	868/898 (97%)	857 (99%)	11 (1%)	61	82
All	All	3471/3592 (97%)	3423 (99%)	48 (1%)	59	81

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	GLU
1	A	267	VAL
1	A	319	ASP
1	A	519	SER
1	A	546	LEU
1	A	604	ASN
1	A	672	VAL
1	A	735	HIS
1	A	737	ILE
1	A	773	LYS
1	A	797	GLU
1	A	956	GLN
1	B	44	THR
1	B	80	GLU
1	B	478	VAL
1	B	519	SER
1	B	546	LEU
1	B	672	VAL
1	B	737	ILE
1	B	850	PHE
1	B	923	SER
1	B	937	LEU
1	B	956	GLN
1	B	1004	SER
1	C	75	GLU
1	C	80	GLU
1	C	131	GLU
1	C	189	LEU
1	C	519	SER
1	C	651	LEU
1	C	672	VAL
1	C	737	ILE
1	C	748	CYS
1	C	751	LEU
1	C	773	LYS
1	C	850	PHE
1	C	956	GLN
1	D	80	GLU
1	D	478	VAL
1	D	519	SER
1	D	546	LEU
1	D	672	VAL
1	D	737	ILE

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Mol	Chain	Res	Type
1	D	744	GLU
1	D	773	LYS
1	D	916	ASP
1	D	956	GLN
1	D	965	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	93	HIS
1	A	128	ASN
1	A	221	GLN
1	A	363	HIS
1	A	394	ASN
1	A	485	GLN
1	A	623	GLN
1	A	713	HIS
1	A	775	GLN
1	A	843	GLN
1	A	863	GLN
1	A	974	HIS
1	B	128	ASN
1	B	135	GLN
1	B	628	GLN
1	B	713	HIS
1	B	739	HIS
1	B	956	GLN
1	C	226	HIS
1	C	445	GLN
1	C	485	GLN
1	C	713	HIS
1	C	885	ASN
1	C	903	GLN
1	C	945	ASN
1	C	950	GLN
1	C	974	HIS
1	D	50	GLN
1	D	93	HIS
1	D	128	ASN
1	D	135	GLN
1	D	221	GLN

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Mol	Chain	Res	Type
1	D	485	GLN
1	D	634	GLN
1	D	693	GLN
1	D	702	GLN
1	D	878	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 85 ligands modelled in this entry, 18 are monoatomic - leaving 67 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	8003	-	3,3,3	2.68	1 (33%)	3,3,3	0.45	0
5	DMS	B	8006	-	3,3,3	2.69	1 (33%)	3,3,3	0.52	0
5	DMS	D	8002	-	3,3,3	2.65	1 (33%)	3,3,3	0.61	0
5	DMS	D	8004	-	3,3,3	2.70	1 (33%)	3,3,3	0.62	0
5	DMS	A	8006	-	3,3,3	2.69	1 (33%)	3,3,3	0.60	0
5	DMS	A	8002	-	3,3,3	2.67	1 (33%)	3,3,3	0.63	0
5	DMS	D	8013	-	3,3,3	2.72	1 (33%)	3,3,3	0.58	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DMS	D	8009	-	3,3,3	2.75	1 (33%)	3,3,3	0.58	0
5	DMS	A	8016	-	3,3,3	2.76	1 (33%)	3,3,3	0.67	0
5	DMS	B	8010	-	3,3,3	2.72	1 (33%)	3,3,3	0.65	0
5	DMS	C	8011	-	3,3,3	2.70	1 (33%)	3,3,3	0.59	0
5	DMS	A	8013	-	3,3,3	2.70	1 (33%)	3,3,3	0.59	0
2	IPT	D	2001	4	15,15,15	0.74	0	18,21,21	0.64	0
5	DMS	C	8010	-	3,3,3	2.77	1 (33%)	3,3,3	0.58	0
5	DMS	B	8002	-	3,3,3	2.68	1 (33%)	3,3,3	0.70	0
5	DMS	B	8004	-	3,3,3	2.72	1 (33%)	3,3,3	0.58	0
5	DMS	A	8005	-	3,3,3	2.63	1 (33%)	3,3,3	0.61	0
5	DMS	B	8009	-	3,3,3	2.77	1 (33%)	3,3,3	0.61	0
5	DMS	D	8012	-	3,3,3	2.72	1 (33%)	3,3,3	0.51	0
5	DMS	D	8014	-	3,3,3	2.75	1 (33%)	3,3,3	0.63	0
5	DMS	B	8007	-	3,3,3	2.76	1 (33%)	3,3,3	0.55	0
5	DMS	C	8012	-	3,3,3	2.71	1 (33%)	3,3,3	0.69	0
2	IPT	B	2001	4	15,15,15	0.70	0	18,21,21	0.65	0
2	IPT	C	2001	4	15,15,15	0.79	1 (6%)	18,21,21	0.70	0
5	DMS	D	8003	-	3,3,3	2.71	1 (33%)	3,3,3	0.49	0
5	DMS	C	8007	-	3,3,3	2.72	1 (33%)	3,3,3	0.65	0
5	DMS	C	8018	-	3,3,3	2.71	1 (33%)	3,3,3	0.56	0
5	DMS	D	8011	-	3,3,3	2.65	1 (33%)	3,3,3	0.53	0
5	DMS	D	8010	-	3,3,3	2.70	1 (33%)	3,3,3	0.59	0
5	DMS	B	8012	-	3,3,3	2.69	1 (33%)	3,3,3	0.52	0
5	DMS	A	8009	-	3,3,3	2.71	1 (33%)	3,3,3	0.55	0
2	IPT	A	2001	4	15,15,15	0.74	0	18,21,21	0.70	0
5	DMS	C	8003	-	3,3,3	2.64	1 (33%)	3,3,3	0.46	0
5	DMS	C	8001	-	3,3,3	2.68	1 (33%)	3,3,3	0.68	0
5	DMS	B	8005	-	3,3,3	2.68	1 (33%)	3,3,3	0.52	0
5	DMS	C	8002	-	3,3,3	2.69	1 (33%)	3,3,3	0.69	0
5	DMS	A	8008	-	3,3,3	2.72	1 (33%)	3,3,3	0.57	0
5	DMS	B	8011	-	3,3,3	2.64	1 (33%)	3,3,3	0.55	0
5	DMS	C	8006	-	3,3,3	2.70	1 (33%)	3,3,3	0.56	0
5	DMS	C	8004	-	3,3,3	2.67	1 (33%)	3,3,3	0.53	0
5	DMS	C	8005	-	3,3,3	2.65	1 (33%)	3,3,3	0.64	0
5	DMS	A	8012	-	3,3,3	2.76	1 (33%)	3,3,3	0.71	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DMS	C	8016	-	3,3,3	2.76	1 (33%)	3,3,3	0.60	0
5	DMS	D	8007	-	3,3,3	2.69	1 (33%)	3,3,3	0.60	0
5	DMS	D	8017	-	3,3,3	2.72	1 (33%)	3,3,3	0.53	0
5	DMS	C	8013	-	3,3,3	2.69	1 (33%)	3,3,3	0.53	0
5	DMS	D	8008	-	3,3,3	2.67	1 (33%)	3,3,3	0.69	0
5	DMS	C	8009	-	3,3,3	2.73	1 (33%)	3,3,3	0.73	0
5	DMS	A	8007	-	3,3,3	2.70	1 (33%)	3,3,3	0.59	0
5	DMS	B	8001	-	3,3,3	2.70	1 (33%)	3,3,3	0.67	0
5	DMS	C	8017	-	3,3,3	2.69	1 (33%)	3,3,3	0.33	0
5	DMS	D	8015	-	3,3,3	2.73	1 (33%)	3,3,3	0.61	0
5	DMS	B	8008	-	3,3,3	2.71	1 (33%)	3,3,3	0.66	0
5	DMS	C	8015	-	3,3,3	2.72	1 (33%)	3,3,3	0.55	0
5	DMS	A	8011	-	3,3,3	2.72	1 (33%)	3,3,3	0.57	0
5	DMS	B	8013	-	3,3,3	2.72	1 (33%)	3,3,3	0.53	0
5	DMS	A	8003	-	3,3,3	2.69	1 (33%)	3,3,3	0.46	0
5	DMS	C	8014	-	3,3,3	2.74	1 (33%)	3,3,3	0.66	0
5	DMS	D	8006	-	3,3,3	2.77	1 (33%)	3,3,3	0.66	0
5	DMS	B	8015	-	3,3,3	2.74	1 (33%)	3,3,3	0.55	0
5	DMS	A	8004	-	3,3,3	2.72	1 (33%)	3,3,3	0.52	0
5	DMS	D	8005	-	3,3,3	2.72	1 (33%)	3,3,3	0.45	0
5	DMS	B	8014	-	3,3,3	2.76	1 (33%)	3,3,3	0.53	0
5	DMS	A	8001	-	3,3,3	2.69	1 (33%)	3,3,3	0.84	0
5	DMS	C	8008	-	3,3,3	2.76	1 (33%)	3,3,3	0.69	0
5	DMS	A	8010	-	3,3,3	2.69	1 (33%)	3,3,3	0.51	0
5	DMS	D	8001	-	3,3,3	2.69	1 (33%)	3,3,3	0.75	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IPT	A	2001	4	-	1/6/26/26	0/1/1/1
2	IPT	B	2001	4	-	2/6/26/26	0/1/1/1
2	IPT	C	2001	4	-	1/6/26/26	0/1/1/1
2	IPT	D	2001	4	-	1/6/26/26	0/1/1/1

All (64) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	8007	DMS	O-S	4.64	1.80	1.50
5	D	8006	DMS	O-S	4.64	1.80	1.50
5	B	8014	DMS	O-S	4.64	1.80	1.50
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	C	8008	DMS	O-S	4.63	1.80	1.50
5	B	8009	DMS	O-S	4.63	1.80	1.50
5	A	8012	DMS	O-S	4.61	1.80	1.50
5	A	8016	DMS	O-S	4.61	1.80	1.50
5	B	8015	DMS	O-S	4.61	1.80	1.50
5	D	8009	DMS	O-S	4.61	1.80	1.50
5	D	8014	DMS	O-S	4.60	1.80	1.50
5	C	8014	DMS	O-S	4.59	1.80	1.50
5	D	8015	DMS	O-S	4.58	1.80	1.50
5	C	8009	DMS	O-S	4.58	1.80	1.50
5	B	8004	DMS	O-S	4.57	1.80	1.50
5	D	8005	DMS	O-S	4.57	1.80	1.50
5	C	8007	DMS	O-S	4.56	1.80	1.50
5	D	8013	DMS	O-S	4.56	1.80	1.50
5	D	8003	DMS	O-S	4.56	1.80	1.50
5	D	8017	DMS	O-S	4.56	1.80	1.50
5	B	8013	DMS	O-S	4.56	1.80	1.50
5	A	8004	DMS	O-S	4.56	1.80	1.50
5	D	8012	DMS	O-S	4.55	1.80	1.50
5	C	8015	DMS	O-S	4.55	1.80	1.50
5	C	8006	DMS	O-S	4.55	1.80	1.50
5	B	8001	DMS	O-S	4.55	1.80	1.50
5	A	8008	DMS	O-S	4.55	1.80	1.50
5	A	8011	DMS	O-S	4.54	1.80	1.50
5	C	8012	DMS	O-S	4.54	1.80	1.50
5	C	8018	DMS	O-S	4.54	1.80	1.50
5	A	8001	DMS	O-S	4.54	1.80	1.50
5	D	8004	DMS	O-S	4.54	1.80	1.50
5	B	8010	DMS	O-S	4.54	1.80	1.50
5	C	8017	DMS	O-S	4.54	1.80	1.50
5	C	8002	DMS	O-S	4.53	1.80	1.50
5	A	8009	DMS	O-S	4.53	1.80	1.50
5	A	8003	DMS	O-S	4.53	1.80	1.50
5	B	8008	DMS	O-S	4.53	1.80	1.50
5	B	8006	DMS	O-S	4.52	1.80	1.50
5	A	8010	DMS	O-S	4.52	1.80	1.50
5	C	8011	DMS	O-S	4.52	1.80	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	8001	DMS	O-S	4.52	1.80	1.50
5	A	8006	DMS	O-S	4.51	1.80	1.50
5	B	8012	DMS	O-S	4.51	1.80	1.50
5	B	8003	DMS	O-S	4.51	1.80	1.50
5	A	8007	DMS	O-S	4.51	1.80	1.50
5	D	8007	DMS	O-S	4.51	1.80	1.50
5	A	8002	DMS	O-S	4.51	1.80	1.50
5	D	8010	DMS	O-S	4.51	1.80	1.50
5	C	8013	DMS	O-S	4.50	1.80	1.50
5	A	8013	DMS	O-S	4.50	1.79	1.50
5	B	8002	DMS	O-S	4.50	1.79	1.50
5	C	8004	DMS	O-S	4.49	1.79	1.50
5	C	8001	DMS	O-S	4.49	1.79	1.50
5	B	8005	DMS	O-S	4.48	1.79	1.50
5	D	8002	DMS	O-S	4.47	1.79	1.50
5	D	8008	DMS	O-S	4.46	1.79	1.50
5	C	8003	DMS	O-S	4.45	1.79	1.50
5	C	8005	DMS	O-S	4.43	1.79	1.50
5	D	8011	DMS	O-S	4.43	1.79	1.50
5	B	8011	DMS	O-S	4.42	1.79	1.50
5	A	8005	DMS	O-S	4.40	1.79	1.50
2	C	2001	IPT	C1-S1	-2.17	1.76	1.80

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	2001	IPT	O5-C5-C6-O6
2	B	2001	IPT	C4-C5-C6-O6
2	A	2001	IPT	O5-C5-C6-O6
2	C	2001	IPT	O5-C5-C6-O6
2	D	2001	IPT	O5-C5-C6-O6

There are no ring outliers.

15 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	8002	DMS	1	0
5	D	8013	DMS	2	0
5	A	8016	DMS	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	8007	DMS	1	0
5	C	8012	DMS	1	0
5	B	8012	DMS	2	0
5	C	8001	DMS	2	0
5	C	8002	DMS	1	0
5	C	8004	DMS	1	0
5	C	8017	DMS	1	0
5	C	8015	DMS	1	0
5	A	8003	DMS	2	0
5	D	8005	DMS	1	0
5	A	8001	DMS	2	0
5	D	8001	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	1014/1052 (96%)	-0.11	22 (2%) 62 57	27, 49, 81, 114	0
1	B	1015/1052 (96%)	-0.04	19 (1%) 66 61	30, 51, 82, 107	0
1	C	1015/1052 (96%)	-0.07	26 (2%) 57 51	29, 51, 84, 111	0
1	D	1015/1052 (96%)	-0.19	22 (2%) 62 57	27, 47, 81, 116	0
All	All	4059/4208 (96%)	-0.10	89 (2%) 62 57	27, 50, 82, 116	0

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	11	LEU	5.1
1	A	10	VAL	5.0
1	B	10	VAL	3.9
1	C	11	LEU	3.8
1	D	733	ALA	3.7
1	D	10	VAL	3.6
1	C	730	LEU	3.5
1	B	798	ALA	3.4
1	B	745	MET	3.3
1	A	730	LEU	3.2
1	D	731	PRO	3.2
1	C	743	SER	3.2
1	D	686	PRO	3.1
1	B	731	PRO	3.1
1	A	11	LEU	3.0
1	B	730	LEU	3.0
1	B	686	PRO	3.0
1	C	736	ALA	3.0
1	C	1023	LYS	2.9
1	A	774	LYS	2.9
1	C	1022	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	737	ILE	2.7
1	D	730	LEU	2.7
1	D	735	HIS	2.7
1	C	731	PRO	2.7
1	B	9	VAL	2.7
1	A	735	HIS	2.7
1	C	10	VAL	2.7
1	D	739	HIS	2.7
1	C	745	MET	2.6
1	D	9	VAL	2.6
1	C	595	THR	2.6
1	C	729	THR	2.6
1	D	76	CYS	2.6
1	A	686	PRO	2.6
1	B	179	ALA	2.6
1	A	801	ILE	2.5
1	C	734	SER	2.5
1	A	773	LYS	2.5
1	D	11	LEU	2.5
1	D	736	ALA	2.5
1	D	732	ALA	2.4
1	D	831	ALA	2.4
1	B	739	HIS	2.4
1	B	743	SER	2.4
1	D	655	MET	2.4
1	B	1022	GLN	2.4
1	C	9	VAL	2.4
1	A	733	ALA	2.4
1	D	1022	GLN	2.4
1	B	733	ALA	2.4
1	B	737	ILE	2.3
1	B	799	THR	2.3
1	A	739	HIS	2.3
1	C	128	ASN	2.3
1	D	1023	LYS	2.3
1	A	12	GLN	2.3
1	D	771	GLY	2.3
1	C	799	THR	2.3
1	A	736	ALA	2.2
1	A	798	ALA	2.2
1	C	733	ALA	2.2
1	C	977	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	596	PRO	2.2
1	C	135	GLN	2.2
1	D	798	ALA	2.2
1	A	76	CYS	2.2
1	A	731	PRO	2.2
1	B	1023	LYS	2.2
1	D	737	ILE	2.1
1	A	931	PHE	2.1
1	B	732	ALA	2.1
1	A	71	GLU	2.1
1	D	699	ARG	2.1
1	B	729	THR	2.1
1	C	773	LYS	2.1
1	A	772	ASP	2.1
1	D	829	THR	2.1
1	C	74	LEU	2.1
1	A	77	ASP	2.0
1	C	131	GLU	2.0
1	A	655	MET	2.0
1	C	76	CYS	2.0
1	B	755	ARG	2.0
1	C	798	ALA	2.0
1	D	734	SER	2.0
1	C	686	PRO	2.0
1	A	685	LEU	2.0
1	C	753	ASN	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	A	8011	4/4	0.70	0.33	108,109,110,111	0
5	DMS	C	8010	4/4	0.72	0.36	114,115,115,117	0
5	DMS	B	8008	4/4	0.74	0.28	87,90,92,92	0
5	DMS	D	8009	4/4	0.75	0.31	116,117,117,117	0
5	DMS	C	8018	4/4	0.79	0.29	98,98,100,101	0
5	DMS	C	8014	4/4	0.79	0.30	90,91,93,96	0
5	DMS	A	8012	4/4	0.80	0.31	97,98,99,99	0
5	DMS	A	8010	4/4	0.81	0.28	81,84,86,86	0
5	DMS	A	8008	4/4	0.81	0.27	93,94,96,96	0
5	DMS	B	8010	4/4	0.82	0.26	85,87,87,89	0
5	DMS	B	8013	4/4	0.83	0.20	112,112,113,113	0
5	DMS	D	8014	4/4	0.83	0.27	94,95,96,98	0
5	DMS	B	8007	4/4	0.84	0.26	97,100,100,100	0
4	NA	A	3103	1/1	0.84	0.20	72,72,72,72	0
5	DMS	C	8016	4/4	0.84	0.23	96,96,97,99	0
5	DMS	C	8007	4/4	0.85	0.24	93,94,95,98	0
5	DMS	D	8006	4/4	0.85	0.21	85,86,88,90	0
5	DMS	C	8015	4/4	0.85	0.28	82,84,85,88	0
5	DMS	C	8011	4/4	0.85	0.24	92,93,93,94	0
5	DMS	A	8013	4/4	0.86	0.25	93,95,96,96	0
5	DMS	D	8012	4/4	0.87	0.25	106,107,108,108	0
5	DMS	D	8013	4/4	0.87	0.29	111,112,113,115	0
5	DMS	A	8007	4/4	0.87	0.21	74,75,76,78	0
5	DMS	A	8016	4/4	0.88	0.20	75,75,75,78	0
5	DMS	C	8013	4/4	0.88	0.22	93,93,94,95	0
5	DMS	D	8010	4/4	0.88	0.23	87,87,88,89	0
5	DMS	D	8008	4/4	0.89	0.22	77,79,79,79	0
5	DMS	C	8008	4/4	0.90	0.17	72,73,74,76	0
5	DMS	B	8005	4/4	0.90	0.19	67,70,71,72	0
5	DMS	A	8004	4/4	0.90	0.18	76,77,79,81	0
5	DMS	C	8006	4/4	0.90	0.25	110,111,112,114	0
3	MG	B	3002	1/1	0.90	0.14	57,57,57,57	0
5	DMS	C	8004	4/4	0.91	0.22	86,88,90,91	0
5	DMS	C	8017	4/4	0.91	0.17	68,74,74,74	0
5	DMS	B	8012	4/4	0.92	0.19	74,78,79,82	0
5	DMS	B	8004	4/4	0.92	0.20	80,81,82,83	0
5	DMS	C	8009	4/4	0.92	0.15	71,71,73,77	0
2	IPT	B	2001	15/15	0.93	0.12	59,70,72,75	0
5	DMS	A	8009	4/4	0.93	0.16	73,74,76,77	0
5	DMS	B	8014	4/4	0.93	0.15	73,75,76,78	0
5	DMS	C	8003	4/4	0.93	0.20	70,71,75,75	0
5	DMS	B	8009	4/4	0.93	0.14	61,62,63,69	0
4	NA	C	3104	1/1	0.93	0.08	57,57,57,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	DMS	D	8015	4/4	0.93	0.15	66,69,72,74	0
5	DMS	D	8004	4/4	0.94	0.17	68,68,71,71	0
5	DMS	B	8006	4/4	0.94	0.16	100,101,102,102	0
5	DMS	D	8007	4/4	0.94	0.17	80,80,80,81	0
5	DMS	C	8002	4/4	0.94	0.14	53,61,62,64	0
2	IPT	C	2001	15/15	0.94	0.12	52,66,71,75	0
5	DMS	A	8005	4/4	0.94	0.14	53,53,55,59	0
5	DMS	D	8011	4/4	0.94	0.14	61,63,66,66	0
5	DMS	B	8002	4/4	0.94	0.16	63,65,67,68	0
5	DMS	B	8003	4/4	0.94	0.15	64,64,66,67	0
2	IPT	A	2001	15/15	0.94	0.12	54,62,72,74	0
5	DMS	A	8002	4/4	0.94	0.16	57,64,65,66	0
5	DMS	D	8017	4/4	0.94	0.17	71,72,72,73	0
4	NA	B	3101	1/1	0.95	0.07	60,60,60,60	0
5	DMS	A	8006	4/4	0.95	0.13	67,68,69,70	0
5	DMS	B	8015	4/4	0.95	0.13	71,72,73,74	0
3	MG	D	3002	1/1	0.95	0.10	53,53,53,53	0
4	NA	A	3101	1/1	0.95	0.10	67,67,67,67	0
3	MG	A	3002	1/1	0.95	0.07	48,48,48,48	0
5	DMS	D	8005	4/4	0.95	0.13	56,57,57,58	0
5	DMS	C	8012	4/4	0.95	0.14	64,64,64,67	0
5	DMS	C	8005	4/4	0.95	0.15	58,59,62,64	0
2	IPT	D	2001	15/15	0.96	0.10	53,60,66,66	0
5	DMS	B	8011	4/4	0.96	0.15	70,72,73,75	0
4	NA	D	3101	1/1	0.96	0.07	58,58,58,58	0
4	NA	B	3102	1/1	0.97	0.07	44,44,44,44	0
4	NA	C	3101	1/1	0.97	0.05	58,58,58,58	0
4	NA	C	3102	1/1	0.97	0.06	42,42,42,42	0
4	NA	A	3102	1/1	0.97	0.05	37,37,37,37	0
5	DMS	D	8002	4/4	0.97	0.13	76,76,77,78	0
5	DMS	B	8001	4/4	0.97	0.10	57,61,63,66	0
3	MG	A	3001	1/1	0.97	0.05	51,51,51,51	0
5	DMS	A	8001	4/4	0.97	0.10	52,52,52,58	0
3	MG	C	3002	1/1	0.97	0.09	58,58,58,58	0
5	DMS	C	8001	4/4	0.98	0.10	60,62,62,64	0
5	DMS	A	8003	4/4	0.98	0.10	68,69,71,71	0
5	DMS	D	8001	4/4	0.98	0.07	42,47,49,50	0
3	MG	D	3001	1/1	0.98	0.04	45,45,45,45	0
5	DMS	D	8003	4/4	0.98	0.12	68,69,70,71	0
4	NA	D	3102	1/1	0.98	0.04	38,38,38,38	0
3	MG	B	3001	1/1	0.99	0.04	57,57,57,57	0
3	MG	C	3001	1/1	0.99	0.04	52,52,52,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.