



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 09:05 PM UTC

PDB ID : 3VD4 / pdb_00003vd4
Title : E. coli (lacZ) beta-galactosidase (N460D) 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 : 2.00 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

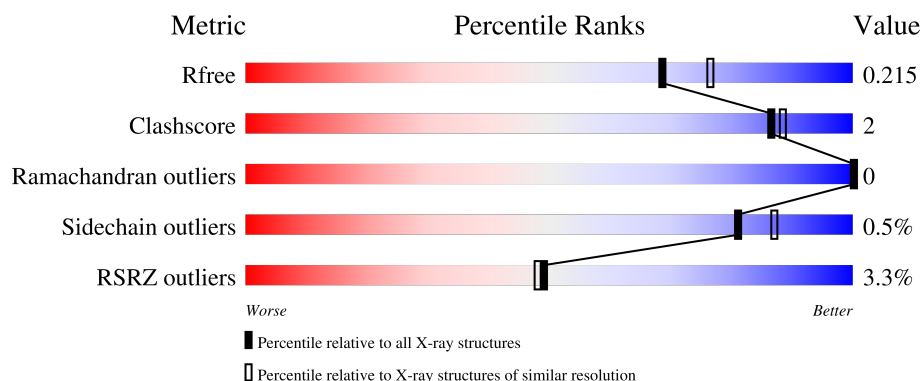
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

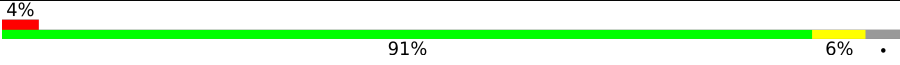
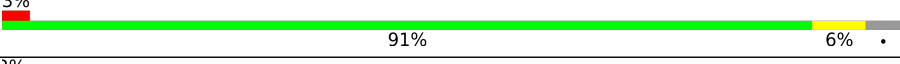
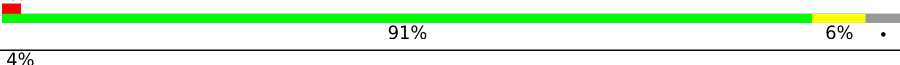
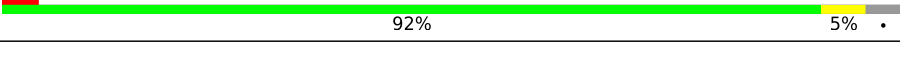
The reported resolution of this entry is 2.00 Å.

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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 36283 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
			8157	5159	1444	1516	38			
1	B	1015	Total	C	N	O	S	0	0	0
			8157	5159	1444	1516	38			
1	C	1015	Total	C	N	O	S	0	0	0
			8157	5159	1444	1516	38			
1	D	1015	Total	C	N	O	S	0	0	0
			8157	5159	1444	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	460	ASP	ASN	engineered mutation	UNP P00722
B	-28	MET	-	expression tag	UNP P00722
B	-27	GLY	-	expression tag	UNP P00722
B	-26	GLY	-	expression tag	UNP P00722
B	-25	SER	-	expression tag	UNP P00722
B	-24	HIS	-	expression tag	UNP P00722
B	-23	HIS	-	expression tag	UNP P00722
B	-22	HIS	-	expression tag	UNP P00722
B	-21	HIS	-	expression tag	UNP P00722
B	-20	HIS	-	expression tag	UNP P00722
B	-19	HIS	-	expression tag	UNP P00722
B	-18	GLY	-	expression tag	UNP P00722
B	-17	MET	-	expression tag	UNP P00722
B	-16	ALA	-	expression tag	UNP P00722
B	-15	SER	-	expression tag	UNP P00722
B	-14	MET	-	expression tag	UNP P00722
B	-13	THR	-	expression tag	UNP P00722
B	-12	GLY	-	expression tag	UNP P00722
B	-11	GLY	-	expression tag	UNP P00722
B	-10	GLN	-	expression tag	UNP P00722
B	-9	GLN	-	expression tag	UNP P00722
B	-8	MET	-	expression tag	UNP P00722
B	-7	GLY	-	expression tag	UNP P00722
B	-6	ARG	-	expression tag	UNP P00722
B	-5	ASP	-	expression tag	UNP P00722
B	-4	LEU	-	expression tag	UNP P00722

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

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

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

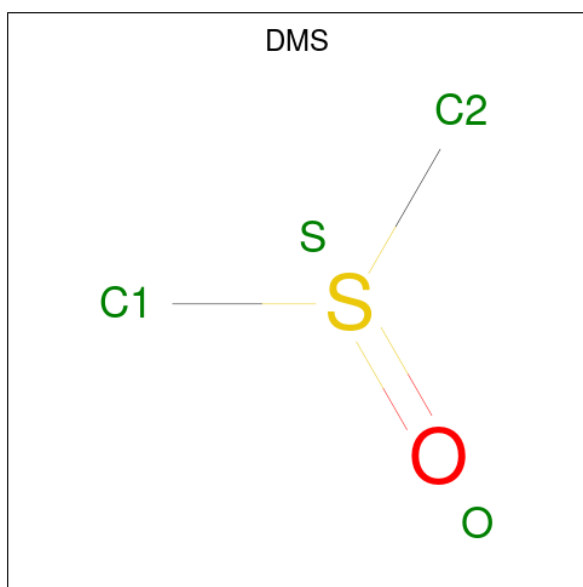
- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

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



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

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

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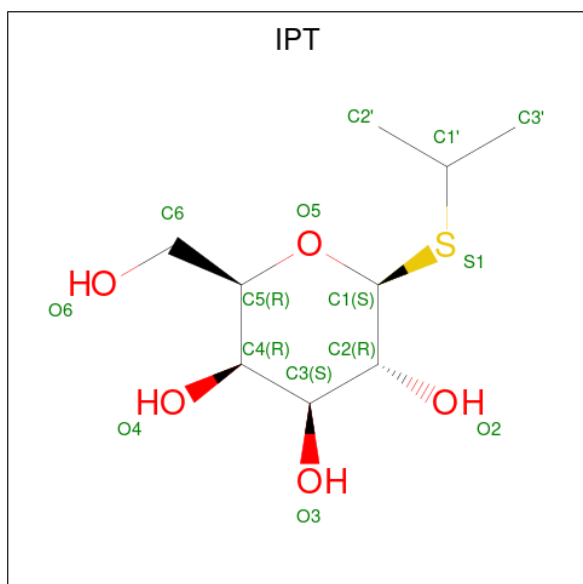
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total 4	C 2	O 1	S 1	0	0
4	C	1	Total 4	C 2	O 1	S 1	0	0
4	C	1	Total 4	C 2	O 1	S 1	0	0
4	C	1	Total 4	C 2	O 1	S 1	0	0
4	C	1	Total 4	C 2	O 1	S 1	0	0
4	C	1	Total 4	C 2	O 1	S 1	0	0
4	C	1	Total 4	C 2	O 1	S 1	0	0
4	C	1	Total 4	C 2	O 1	S 1	0	0
4	C	1	Total 4	C 2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

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

- Molecule 5 is 1-methylethyl 1-thio-beta-D-galactopyranoside (CCD ID: IPT) (formula: $C_9H_{18}O_5S$).



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

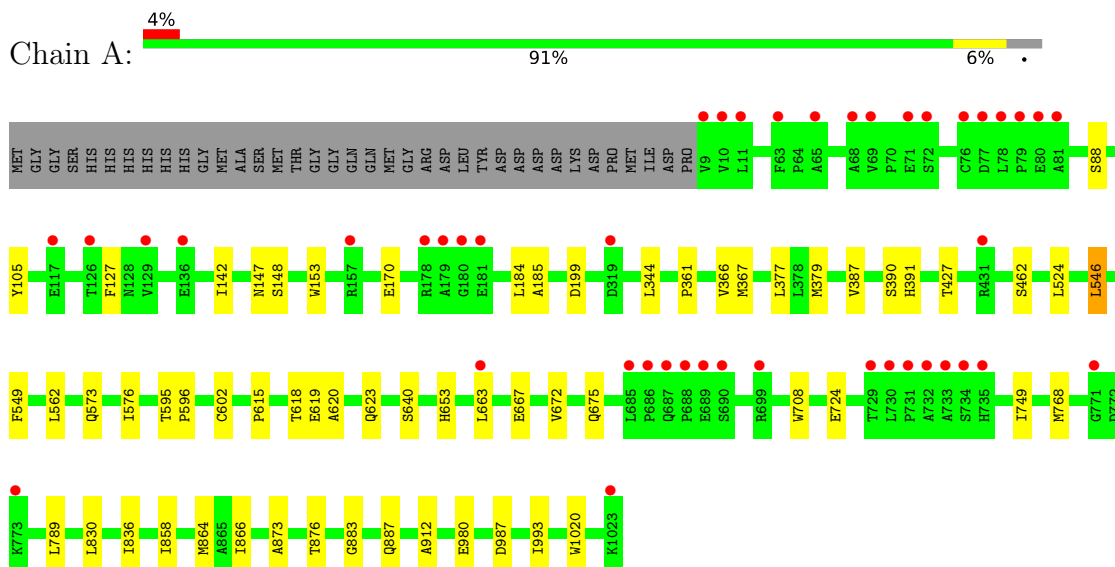
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	750	Total	O	0	0
			750	750		
6	B	909	Total	O	0	0
			909	909		
6	C	897	Total	O	0	0
			897	897		
6	D	744	Total	O	0	0
			744	744		

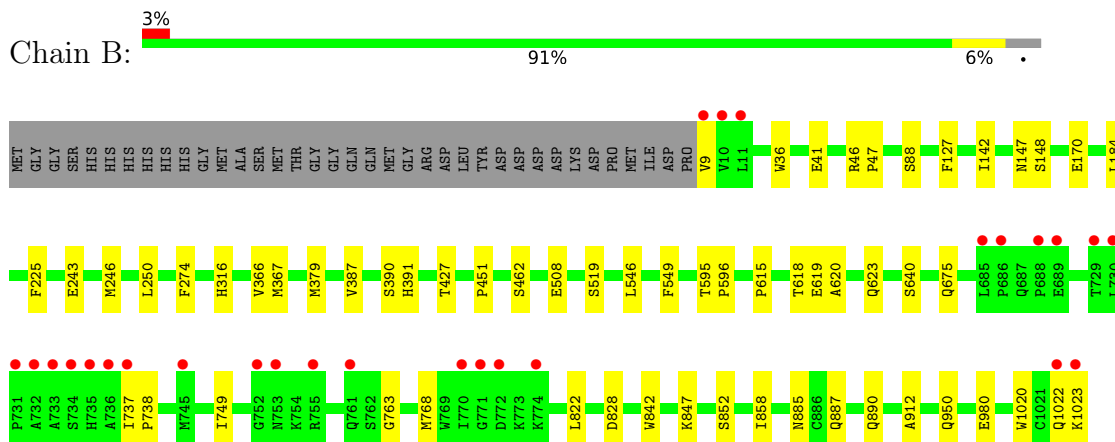
3 Residue-property plots [i](#)

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

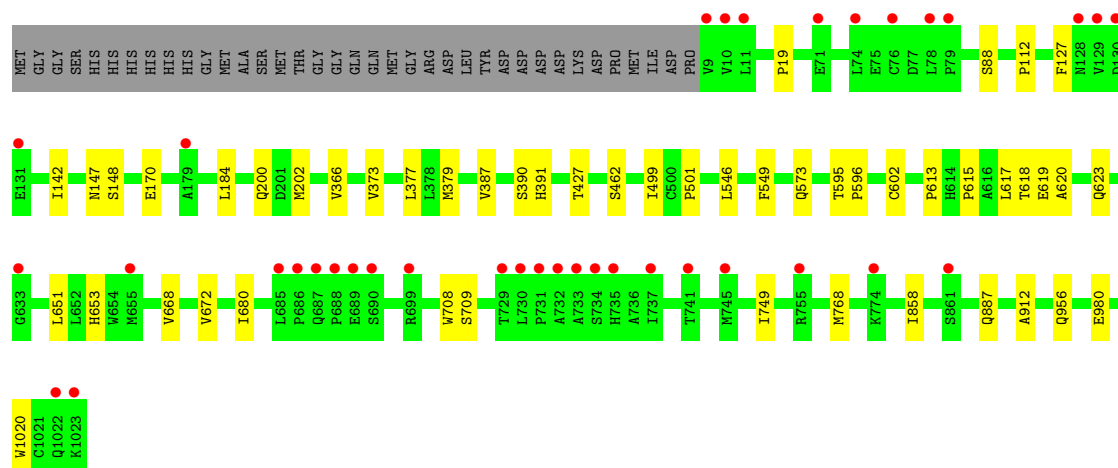


- Molecule 1: Beta-galactosidase



- Molecule 1: Beta-galactosidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	151.76Å 162.45Å 203.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	110.90 – 2.00 110.90 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (110.90-2.00) 92.8 (110.90-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0109, CNS	Depositor
R, R_{free}	0.172 , 0.212 0.175 , 0.215	Depositor DCC
R_{free} test set	4492 reflections (1.44%)	wwPDB-VP
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	36283	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.47 % 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 2.7241e-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: MG, IPT, NA, DMS

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	0.46	0/8399	0.68	0/11458
1	B	0.48	0/8399	0.68	0/11458
1	C	0.48	0/8399	0.68	0/11458
1	D	0.45	0/8399	0.68	0/11458
All	All	0.47	0/33596	0.68	0/45832

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	8157	0	7751	34	0
1	B	8157	0	7751	39	0
1	C	8157	0	7751	34	0
1	D	8157	0	7751	27	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	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	4	0	0	0	0
3	C	4	0	0	0	0
3	D	3	0	0	0	0
4	A	68	0	102	2	0
4	B	68	0	102	8	0
4	C	76	0	114	8	0
4	D	60	0	90	2	0
5	A	15	0	17	0	0
5	B	15	0	17	0	0
5	C	15	0	17	0	0
5	D	15	0	17	0	0
6	A	750	0	0	2	0
6	B	909	0	0	3	0
6	C	897	0	0	0	0
6	D	744	0	0	0	0
All	All	36283	0	31480	133	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 (133) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:379:MET:HE1	1:B:387:VAL:HB	1.75	0.68
1:D:379:MET:HE1	1:D:387:VAL:HB	1.79	0.63
1:C:508:GLU:OE2	4:C:8006:DMS:H11	1.99	0.63
1:A:142:ILE:HG12	1:A:170:GLU:HG2	1.80	0.62
1:C:377:LEU:HD22	1:C:708:TRP:HA	1.80	0.62
1:A:768:MET:HE1	1:A:1020:TRP:CZ2	2.34	0.62
1:D:887:GLN:NE2	1:D:980:GLU:O	2.33	0.62
1:C:749:ILE:HD13	1:C:858:ILE:HD12	1.82	0.60
1:C:887:GLN:NE2	1:C:980:GLU:O	2.33	0.59
1:A:379:MET:HE1	1:A:387:VAL:HB	1.84	0.58
1:B:749:ILE:HD13	1:B:858:ILE:HD12	1.87	0.57
1:B:887:GLN:NE2	1:B:980:GLU:O	2.35	0.57
1:B:508:GLU:OE2	4:B:8006:DMS:H21	2.05	0.56
1:D:615:PRO:O	1:D:618:THR:HG22	2.06	0.56
4:C:8003:DMS:O	4:C:8015:DMS:H12	2.08	0.54
1:A:887:GLN:NE2	1:A:980:GLU:O	2.39	0.54
1:B:950:GLN:HB2	1:B:1023:LYS:HB3	1.90	0.53
1:B:47:PRO:HD3	4:B:8016:DMS:H12	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:724:GLU:O	1:B:847:LYS:NZ	2.42	0.52
1:B:367:MET:HA	1:B:367:MET:HE2	1.91	0.52
1:C:451:PRO:HA	4:C:8001:DMS:H23	1.90	0.52
1:A:768:MET:HE1	1:A:1020:TRP:CH2	2.45	0.52
1:A:623:GLN:HA	4:A:8002:DMS:O	2.09	0.51
1:B:46:ARG:NH1	6:B:4794:HOH:O	2.25	0.51
1:B:41:GLU:HG2	1:B:46:ARG:NH1	2.26	0.51
1:A:88:SER:HA	1:A:366:VAL:HG21	1.92	0.50
4:B:8012:DMS:H13	1:C:478:VAL:HG22	1.94	0.50
1:A:427:THR:HG21	1:A:462:SER:HB3	1.92	0.50
1:B:127:PHE:HE1	1:B:184:LEU:HG	1.77	0.49
1:D:708:TRP:CE3	1:D:709:SER:HB3	2.46	0.49
1:A:595:THR:HA	1:A:596:PRO:C	2.37	0.49
1:D:708:TRP:CZ2	4:D:8003:DMS:H23	2.47	0.49
1:C:93:HIS:CG	4:C:8011:DMS:H12	2.48	0.49
1:C:699:ARG:HH12	4:C:8012:DMS:H12	1.78	0.49
1:D:127:PHE:CE1	1:D:184:LEU:HG	2.48	0.49
1:C:127:PHE:CE1	1:C:184:LEU:HG	2.47	0.49
1:C:379:MET:HE1	1:C:387:VAL:HB	1.94	0.49
1:A:619:GLU:HA	1:A:912:ALA:HB2	1.95	0.49
1:A:749:ILE:HD11	1:A:836:ILE:HD11	1.95	0.48
1:C:623:GLN:HA	4:C:8002:DMS:O	2.12	0.48
1:D:499:ILE:HG22	1:D:501:PRO:HD3	1.94	0.48
1:B:451:PRO:HA	4:B:8001:DMS:H23	1.95	0.48
1:B:763:GLY:HA3	1:B:822:LEU:HD13	1.96	0.48
1:B:127:PHE:CE1	1:B:184:LEU:HG	2.48	0.48
1:D:768:MET:HE1	1:D:1020:TRP:CZ2	2.48	0.48
1:B:615:PRO:O	1:B:618:THR:HG22	2.14	0.48
1:D:373:VAL:O	1:D:377:LEU:HG	2.14	0.48
1:D:88:SER:HA	1:D:366:VAL:HG21	1.95	0.47
4:A:8003:DMS:H13	6:A:4073:HOH:O	2.14	0.47
1:D:623:GLN:HA	4:D:8002:DMS:O	2.14	0.47
1:C:367:MET:HE2	1:C:367:MET:HA	1.97	0.47
1:C:950:GLN:HB2	1:C:1023:LYS:HB3	1.97	0.47
1:A:377:LEU:HD22	1:A:708:TRP:HA	1.97	0.46
1:A:640:SER:O	1:A:675:GLN:HA	2.15	0.46
1:B:595:THR:HA	1:B:596:PRO:C	2.39	0.46
1:B:47:PRO:HD3	4:B:8016:DMS:C1	2.46	0.46
1:B:36:TRP:HH2	4:B:8003:DMS:H23	1.81	0.46
1:B:427:THR:HG21	1:B:462:SER:HB3	1.98	0.46
1:A:153:TRP:HB2	1:A:185:ALA:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:PHE:HE1	1:A:184:LEU:HG	1.81	0.46
1:C:615:PRO:O	1:C:618:THR:HG22	2.15	0.46
1:B:88:SER:HA	1:B:366:VAL:HG21	1.98	0.45
1:B:768:MET:HE1	1:B:1020:TRP:CZ2	2.51	0.45
1:A:749:ILE:HD13	1:A:858:ILE:HD12	1.98	0.45
1:B:46:ARG:NH1	6:B:4786:HOH:O	2.47	0.45
1:B:246:MET:HB3	1:B:274:PHE:CZ	2.52	0.45
1:B:1022:GLN:HG2	6:B:4806:HOH:O	2.17	0.45
1:B:246:MET:SD	1:B:250:LEU:HD23	2.57	0.44
1:D:427:THR:HG21	1:D:462:SER:HB3	1.98	0.44
1:C:788:PRO:HD2	1:C:968:MET:HB2	2.00	0.44
1:D:377:LEU:HD22	1:D:708:TRP:HA	1.99	0.44
1:D:573:GLN:HB2	1:D:602:CYS:O	2.17	0.44
1:D:19:PRO:HD3	1:D:112:PRO:HB3	1.99	0.44
1:A:549:PHE:CE2	1:A:620:ALA:HA	2.53	0.44
1:A:830:LEU:HD22	1:B:828:ASP:HB3	2.00	0.44
1:B:9:VAL:HG22	1:C:9:VAL:HG22	2.00	0.44
1:C:595:THR:HA	1:C:596:PRO:C	2.42	0.44
1:B:142:ILE:HG12	1:B:170:GLU:HG2	1.99	0.44
1:D:147:ASN:HA	1:D:148:SER:HA	1.66	0.43
1:C:127:PHE:HE1	1:C:184:LEU:HG	1.82	0.43
1:C:640:SER:O	1:C:675:GLN:HA	2.18	0.43
1:A:546:LEU:HA	6:A:4124:HOH:O	2.17	0.43
1:A:873:ALA:O	1:A:876:THR:HG22	2.18	0.43
1:A:615:PRO:O	1:A:618:THR:HG22	2.19	0.43
1:C:573:GLN:HB2	1:C:602:CYS:O	2.19	0.43
1:D:613:PRO:HB3	1:D:617:LEU:HD23	2.00	0.43
1:C:524:LEU:HD11	1:C:562:LEU:HG	2.00	0.43
1:D:390:SER:HA	1:D:391:HIS:HA	1.83	0.43
1:D:127:PHE:HE1	1:D:184:LEU:HG	1.83	0.43
1:D:668:VAL:HG11	1:D:680:ILE:HG21	1.99	0.43
1:A:573:GLN:HB2	1:A:602:CYS:O	2.19	0.43
1:C:88:SER:HA	1:C:366:VAL:HG21	2.00	0.43
1:C:147:ASN:HA	1:C:148:SER:HA	1.71	0.43
1:C:499:ILE:HG22	1:C:501:PRO:HD3	2.00	0.43
1:C:619:GLU:HA	1:C:912:ALA:HB2	2.01	0.43
1:A:864:MET:HE3	1:A:866:ILE:HD11	2.01	0.42
1:D:749:ILE:HD13	1:D:858:ILE:HD12	2.00	0.42
1:B:623:GLN:NE2	4:B:8002:DMS:H12	2.34	0.42
1:B:316:HIS:CD2	4:B:8005:DMS:H23	2.54	0.42
1:B:749:ILE:CD1	1:B:858:ILE:HD12	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:653:HIS:CD2	1:A:667:GLU:HG2	2.54	0.42
1:D:549:PHE:CE2	1:D:620:ALA:HA	2.54	0.42
1:D:595:THR:HA	1:D:596:PRO:C	2.44	0.42
1:B:619:GLU:HA	1:B:912:ALA:HB2	2.01	0.42
1:C:549:PHE:CE2	1:C:620:ALA:HA	2.54	0.42
1:D:200:GLN:HB2	1:D:202:MET:HE2	2.02	0.42
1:B:147:ASN:HA	1:B:148:SER:HA	1.66	0.41
1:B:549:PHE:CE2	1:B:620:ALA:HA	2.55	0.41
1:D:619:GLU:HA	1:D:912:ALA:HB2	2.01	0.41
1:A:105:TYR:CE1	1:A:199:ASP:HB2	2.55	0.41
1:A:361:PRO:HB2	1:A:576:ILE:HG12	2.02	0.41
1:A:524:LEU:HD11	1:A:562:LEU:HG	2.02	0.41
1:A:147:ASN:HA	1:A:148:SER:HA	1.63	0.41
1:A:367:MET:HA	1:A:367:MET:HE2	2.02	0.41
1:A:789:LEU:HD11	1:A:993:ILE:HG22	2.02	0.41
1:A:127:PHE:CE1	1:A:184:LEU:HG	2.55	0.41
1:B:390:SER:HA	1:B:391:HIS:HA	1.81	0.41
1:A:883:GLY:HA3	1:A:987:ASP:HA	2.01	0.41
1:B:737:ILE:HA	1:B:738:PRO:HD3	1.93	0.41
1:C:143:PHE:O	1:C:168:PRO:HA	2.21	0.41
1:D:142:ILE:HG23	1:D:170:GLU:HG2	2.02	0.41
1:C:84:VAL:HG12	4:C:8011:DMS:H13	2.03	0.41
1:C:907:PRO:HG2	1:C:990:HIS:O	2.20	0.41
1:C:270:GLY:HA3	4:C:8004:DMS:O	2.21	0.41
1:C:749:ILE:CD1	1:C:858:ILE:HD12	2.48	0.41
1:C:948:PRO:O	1:C:1022:GLN:HG3	2.21	0.41
1:B:640:SER:O	1:B:675:GLN:HA	2.21	0.41
1:D:651:LEU:CD1	1:D:653:HIS:HD2	2.34	0.40
1:C:768:MET:HE1	1:C:1020:TRP:CZ2	2.56	0.40
1:B:842:TRP:HZ3	1:B:852:SER:HB3	1.85	0.40
1:A:390:SER:HA	1:A:391:HIS:HA	1.79	0.40
1:B:225:PHE:HA	1:B:243:GLU:O	2.22	0.40
1:C:390:SER:HA	1:C:391:HIS:HA	1.83	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	1013/1052 (96%)	986 (97%)	27 (3%)	0	100	100
1	C	1013/1052 (96%)	988 (98%)	25 (2%)	0	100	100
1	D	1013/1052 (96%)	981 (97%)	32 (3%)	0	100	100
All	All	4052/4208 (96%)	3933 (97%)	119 (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%)	864 (100%)	4 (0%)	81	87
1	B	868/898 (97%)	864 (100%)	4 (0%)	81	87
1	C	868/898 (97%)	862 (99%)	6 (1%)	76	82
1	D	868/898 (97%)	865 (100%)	3 (0%)	86	91
All	All	3472/3592 (97%)	3455 (100%)	17 (0%)	81	87

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	344	LEU
1	A	546	LEU

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Mol	Chain	Res	Type
1	A	663	LEU
1	A	672	VAL
1	B	519	SER
1	B	546	LEU
1	B	885	ASN
1	B	890	GLN
1	C	49	GLN
1	C	519	SER
1	C	546	LEU
1	C	672	VAL
1	C	772	ASP
1	C	956	GLN
1	D	546	LEU
1	D	672	VAL
1	D	956	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	128	ASN
1	A	294	ASN
1	A	363	HIS
1	A	485	GLN
1	A	628	GLN
1	A	702	GLN
1	A	843	GLN
1	A	903	GLN
1	A	945	ASN
1	A	950	GLN
1	A	977	HIS
1	B	23	GLN
1	B	50	GLN
1	B	363	HIS
1	B	485	GLN
1	B	653	HIS
1	B	843	GLN
1	B	890	GLN
1	B	903	GLN
1	B	977	HIS
1	B	1017	GLN
1	B	1022	GLN

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Mol	Chain	Res	Type
1	C	49	GLN
1	C	128	ASN
1	C	365	GLN
1	C	381	GLN
1	C	485	GLN
1	C	843	GLN
1	C	1022	GLN
1	D	128	ASN
1	D	297	ASN
1	D	394	ASN
1	D	445	GLN
1	D	485	GLN
1	D	653	HIS
1	D	718	GLN
1	D	863	GLN
1	D	903	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 95 ligands modelled in this entry, 23 are monoatomic - leaving 72 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DMS	D	8006	-	3,3,3	2.72	1 (33%)	3,3,3	0.43	0
4	DMS	C	8014	-	3,3,3	2.71	1 (33%)	3,3,3	0.56	0
4	DMS	D	8011	-	3,3,3	2.70	1 (33%)	3,3,3	0.54	0
4	DMS	B	8010	-	3,3,3	2.66	1 (33%)	3,3,3	0.40	0
4	DMS	C	8015	-	3,3,3	2.73	1 (33%)	3,3,3	0.53	0
4	DMS	C	8018	-	3,3,3	2.71	1 (33%)	3,3,3	0.56	0
4	DMS	D	8007	-	3,3,3	2.70	1 (33%)	3,3,3	0.63	0
4	DMS	A	8005	-	3,3,3	2.66	1 (33%)	3,3,3	0.47	0
4	DMS	A	8004	-	3,3,3	2.69	1 (33%)	3,3,3	0.58	0
4	DMS	B	8012	-	3,3,3	2.67	1 (33%)	3,3,3	0.52	0
4	DMS	C	8007	-	3,3,3	2.69	1 (33%)	3,3,3	0.63	0
4	DMS	A	8008	-	3,3,3	2.75	1 (33%)	3,3,3	0.53	0
4	DMS	C	8010	-	3,3,3	2.70	1 (33%)	3,3,3	0.63	0
4	DMS	B	8003	-	3,3,3	2.65	1 (33%)	3,3,3	0.65	0
4	DMS	B	8009	-	3,3,3	2.29	1 (33%)	3,3,3	0.22	0
4	DMS	C	8005	-	3,3,3	2.69	1 (33%)	3,3,3	0.53	0
4	DMS	C	8019	-	3,3,3	2.74	1 (33%)	3,3,3	0.52	0
4	DMS	D	8004	-	3,3,3	2.67	1 (33%)	3,3,3	0.55	0
4	DMS	D	8005	-	3,3,3	2.61	1 (33%)	3,3,3	0.36	0
4	DMS	A	8001	-	3,3,3	2.27	1 (33%)	3,3,3	0.50	0
4	DMS	B	8006	-	3,3,3	2.72	1 (33%)	3,3,3	0.46	0
4	DMS	D	8002	-	3,3,3	2.60	1 (33%)	3,3,3	0.43	0
4	DMS	B	8014	-	3,3,3	2.63	1 (33%)	3,3,3	0.23	0
4	DMS	B	8013	-	3,3,3	2.70	1 (33%)	3,3,3	0.53	0
5	IPT	C	2001	3	15,15,15	0.72	0	18,21,21	0.71	0
4	DMS	A	8017	-	3,3,3	2.76	1 (33%)	3,3,3	0.62	0
4	DMS	C	8003	-	3,3,3	2.67	1 (33%)	3,3,3	0.54	0
4	DMS	B	8002	-	3,3,3	2.63	1 (33%)	3,3,3	0.83	0
4	DMS	C	8013	-	3,3,3	2.65	1 (33%)	3,3,3	0.52	0
4	DMS	D	8013	-	3,3,3	2.73	1 (33%)	3,3,3	0.63	0
4	DMS	A	8012	-	3,3,3	2.73	1 (33%)	3,3,3	0.59	0
5	IPT	B	2001	3	15,15,15	0.73	0	18,21,21	0.73	0
4	DMS	B	8004	-	3,3,3	2.66	1 (33%)	3,3,3	0.33	0
4	DMS	B	8005	-	3,3,3	2.73	1 (33%)	3,3,3	0.67	0
4	DMS	B	8016	-	3,3,3	2.72	1 (33%)	3,3,3	0.60	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DMS	B	8017	-	3,3,3	2.71	1 (33%)	3,3,3	0.52	0
4	DMS	B	8008	-	3,3,3	2.70	1 (33%)	3,3,3	0.56	0
4	DMS	C	8006	-	3,3,3	2.71	1 (33%)	3,3,3	0.65	0
4	DMS	C	8012	-	3,3,3	2.71	1 (33%)	3,3,3	0.57	0
4	DMS	C	8002	-	3,3,3	2.59	1 (33%)	3,3,3	0.39	0
4	DMS	B	8001	-	3,3,3	2.64	1 (33%)	3,3,3	0.91	0
4	DMS	D	8015	-	3,3,3	2.70	1 (33%)	3,3,3	0.51	0
4	DMS	A	8016	-	3,3,3	2.74	1 (33%)	3,3,3	0.55	0
5	IPT	D	2001	3	15,15,15	0.73	0	18,21,21	0.77	0
4	DMS	D	8012	-	3,3,3	2.74	1 (33%)	3,3,3	0.53	0
4	DMS	D	8014	-	3,3,3	2.77	1 (33%)	3,3,3	0.61	0
4	DMS	C	8004	-	3,3,3	2.61	1 (33%)	3,3,3	0.29	0
4	DMS	C	8016	-	3,3,3	2.71	1 (33%)	3,3,3	0.60	0
4	DMS	A	8002	-	3,3,3	2.58	1 (33%)	3,3,3	0.46	0
4	DMS	C	8008	-	3,3,3	2.69	1 (33%)	3,3,3	0.67	0
4	DMS	D	8010	-	3,3,3	2.65	1 (33%)	3,3,3	0.46	0
4	DMS	D	8003	-	3,3,3	2.36	1 (33%)	3,3,3	0.43	0
4	DMS	D	8008	-	3,3,3	2.74	1 (33%)	3,3,3	0.51	0
4	DMS	A	8010	-	3,3,3	2.66	1 (33%)	3,3,3	0.47	0
4	DMS	C	8001	-	3,3,3	2.69	1 (33%)	3,3,3	0.98	0
4	DMS	B	8011	-	3,3,3	2.74	1 (33%)	3,3,3	0.68	0
4	DMS	C	8009	-	3,3,3	2.69	1 (33%)	3,3,3	0.57	0
4	DMS	A	8003	-	3,3,3	2.65	1 (33%)	3,3,3	0.57	0
4	DMS	C	8017	-	3,3,3	2.69	1 (33%)	3,3,3	0.54	0
4	DMS	A	8013	-	3,3,3	2.75	1 (33%)	3,3,3	0.55	0
4	DMS	A	8007	-	3,3,3	2.72	1 (33%)	3,3,3	0.65	0
4	DMS	A	8011	-	3,3,3	2.69	1 (33%)	3,3,3	0.49	0
5	IPT	A	2001	3	15,15,15	0.71	0	18,21,21	0.63	0
4	DMS	C	8011	-	3,3,3	2.67	1 (33%)	3,3,3	0.71	0
4	DMS	B	8015	-	3,3,3	2.69	1 (33%)	3,3,3	0.65	0
4	DMS	A	8009	-	3,3,3	2.71	1 (33%)	3,3,3	0.62	0
4	DMS	D	8001	-	3,3,3	2.71	1 (33%)	3,3,3	0.94	0
4	DMS	B	8007	-	3,3,3	2.73	1 (33%)	3,3,3	0.77	0
4	DMS	A	8015	-	3,3,3	2.69	1 (33%)	3,3,3	0.64	0
4	DMS	D	8009	-	3,3,3	2.71	1 (33%)	3,3,3	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DMS	A	8006	-	3,3,3	2.70	1 (33%)	3,3,3	0.48	0
4	DMS	A	8014	-	3,3,3	2.71	1 (33%)	3,3,3	0.52	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
5	IPT	D	2001	3	-	1/6/26/26	0/1/1/1
5	IPT	C	2001	3	-	1/6/26/26	0/1/1/1
5	IPT	B	2001	3	-	1/6/26/26	0/1/1/1
5	IPT	A	2001	3	-	1/6/26/26	0/1/1/1

All (68) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	8017	DMS	O-S	4.64	1.80	1.50
4	D	8014	DMS	O-S	4.64	1.80	1.50
4	A	8013	DMS	O-S	4.61	1.80	1.50
4	A	8008	DMS	O-S	4.61	1.80	1.50
4	C	8019	DMS	O-S	4.60	1.80	1.50
4	B	8007	DMS	O-S	4.60	1.80	1.50
4	B	8005	DMS	O-S	4.60	1.80	1.50
4	A	8016	DMS	O-S	4.60	1.80	1.50
4	D	8012	DMS	O-S	4.60	1.80	1.50
4	B	8011	DMS	O-S	4.59	1.80	1.50
4	D	8008	DMS	O-S	4.59	1.80	1.50
4	C	8015	DMS	O-S	4.59	1.80	1.50
4	D	8013	DMS	O-S	4.58	1.80	1.50
4	A	8012	DMS	O-S	4.58	1.80	1.50
4	B	8016	DMS	O-S	4.57	1.80	1.50
4	C	8006	DMS	O-S	4.57	1.80	1.50
4	B	8006	DMS	O-S	4.57	1.80	1.50
4	A	8007	DMS	O-S	4.56	1.80	1.50
4	C	8012	DMS	O-S	4.55	1.80	1.50
4	D	8006	DMS	O-S	4.55	1.80	1.50
4	C	8014	DMS	O-S	4.55	1.80	1.50
4	C	8016	DMS	O-S	4.55	1.80	1.50
4	A	8014	DMS	O-S	4.54	1.80	1.50
4	B	8017	DMS	O-S	4.54	1.80	1.50
4	A	8009	DMS	O-S	4.54	1.80	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	8018	DMS	O-S	4.54	1.80	1.50
4	D	8015	DMS	O-S	4.54	1.80	1.50
4	D	8007	DMS	O-S	4.54	1.80	1.50
4	D	8009	DMS	O-S	4.53	1.80	1.50
4	C	8010	DMS	O-S	4.53	1.80	1.50
4	D	8001	DMS	O-S	4.53	1.80	1.50
4	A	8004	DMS	O-S	4.53	1.80	1.50
4	B	8013	DMS	O-S	4.52	1.80	1.50
4	D	8011	DMS	O-S	4.52	1.80	1.50
4	B	8008	DMS	O-S	4.52	1.80	1.50
4	C	8017	DMS	O-S	4.52	1.80	1.50
4	C	8001	DMS	O-S	4.52	1.80	1.50
4	C	8005	DMS	O-S	4.51	1.80	1.50
4	C	8007	DMS	O-S	4.51	1.80	1.50
4	A	8011	DMS	O-S	4.51	1.80	1.50
4	C	8008	DMS	O-S	4.51	1.80	1.50
4	A	8015	DMS	O-S	4.51	1.80	1.50
4	A	8006	DMS	O-S	4.50	1.79	1.50
4	C	8009	DMS	O-S	4.50	1.79	1.50
4	B	8015	DMS	O-S	4.49	1.79	1.50
4	C	8003	DMS	O-S	4.48	1.79	1.50
4	D	8004	DMS	O-S	4.48	1.79	1.50
4	B	8012	DMS	O-S	4.48	1.79	1.50
4	B	8010	DMS	O-S	4.47	1.79	1.50
4	C	8011	DMS	O-S	4.47	1.79	1.50
4	A	8005	DMS	O-S	4.47	1.79	1.50
4	A	8003	DMS	O-S	4.46	1.79	1.50
4	B	8003	DMS	O-S	4.46	1.79	1.50
4	A	8010	DMS	O-S	4.45	1.79	1.50
4	B	8004	DMS	O-S	4.45	1.79	1.50
4	D	8010	DMS	O-S	4.44	1.79	1.50
4	C	8013	DMS	O-S	4.43	1.79	1.50
4	B	8014	DMS	O-S	4.43	1.79	1.50
4	B	8001	DMS	O-S	4.43	1.79	1.50
4	B	8002	DMS	O-S	4.41	1.79	1.50
4	C	8004	DMS	O-S	4.37	1.79	1.50
4	D	8002	DMS	O-S	4.37	1.79	1.50
4	D	8005	DMS	O-S	4.37	1.79	1.50
4	C	8002	DMS	O-S	4.35	1.79	1.50
4	A	8002	DMS	O-S	4.33	1.78	1.50
4	D	8003	DMS	O-S	3.99	1.76	1.50
4	B	8009	DMS	O-S	3.86	1.75	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	8001	DMS	O-S	3.83	1.75	1.50

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

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

There are no ring outliers.

19 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	8015	DMS	1	0
4	B	8012	DMS	1	0
4	B	8003	DMS	1	0
4	B	8006	DMS	1	0
4	D	8002	DMS	1	0
4	C	8003	DMS	1	0
4	B	8002	DMS	1	0
4	B	8005	DMS	1	0
4	B	8016	DMS	2	0
4	C	8006	DMS	1	0
4	C	8012	DMS	1	0
4	C	8002	DMS	1	0
4	B	8001	DMS	1	0
4	C	8004	DMS	1	0
4	A	8002	DMS	1	0
4	D	8003	DMS	1	0
4	C	8001	DMS	1	0
4	A	8003	DMS	1	0
4	C	8011	DMS	2	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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%)	-0.02	44 (4%)	40	39	11, 24, 48, 78	0
1	B	1015/1052 (96%)	-0.36	27 (2%)	56	55	10, 19, 42, 76	0
1	C	1015/1052 (96%)	-0.41	24 (2%)	59	59	9, 19, 42, 72	0
1	D	1015/1052 (96%)	-0.10	37 (3%)	46	45	13, 24, 47, 83	0
All	All	4060/4208 (96%)	-0.22	132 (3%)	49	48	9, 22, 45, 83	0

All (132) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	10	VAL	6.0
1	B	770	ILE	5.0
1	A	11	LEU	5.0
1	B	11	LEU	4.9
1	C	10	VAL	4.7
1	A	9	VAL	4.7
1	B	685	LEU	4.7
1	A	733	ALA	4.5
1	A	685	LEU	4.4
1	B	10	VAL	4.3
1	B	731	PRO	4.3
1	D	10	VAL	4.2
1	A	730	LEU	4.2
1	C	686	PRO	4.1
1	C	11	LEU	4.1
1	B	1023	LYS	4.0
1	C	1022	GLN	3.9
1	D	730	LEU	3.8
1	D	689	GLU	3.8
1	B	732	ALA	3.8
1	D	685	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	1022	GLN	3.8
1	A	771	GLY	3.7
1	C	9	VAL	3.7
1	D	688	PRO	3.7
1	D	733	ALA	3.7
1	B	752	GLY	3.7
1	B	771	GLY	3.6
1	B	9	VAL	3.6
1	A	686	PRO	3.6
1	A	77	ASP	3.5
1	D	11	LEU	3.5
1	D	686	PRO	3.4
1	D	732	ALA	3.4
1	A	319	ASP	3.3
1	C	1023	LYS	3.3
1	D	9	VAL	3.3
1	A	71	GLU	3.3
1	D	737	ILE	3.3
1	B	730	LEU	3.3
1	C	770	ILE	3.2
1	C	731	PRO	3.2
1	B	755	ARG	3.2
1	B	772	ASP	3.1
1	A	63	PHE	3.1
1	A	732	ALA	3.1
1	D	129	VAL	3.1
1	A	78	LEU	3.1
1	C	753	ASN	3.1
1	B	689	GLU	3.1
1	B	753	ASN	3.0
1	A	773	LYS	3.0
1	C	685	LEU	3.0
1	A	735	HIS	3.0
1	A	689	GLU	3.0
1	B	774	LYS	2.9
1	D	741	THR	2.9
1	A	81	ALA	2.9
1	A	688	PRO	2.9
1	D	76	CYS	2.8
1	A	79	PRO	2.8
1	D	735	HIS	2.8
1	A	731	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	688	PRO	2.8
1	B	737	ILE	2.7
1	D	1022	GLN	2.7
1	D	687	GLN	2.6
1	A	65	ALA	2.6
1	B	733	ALA	2.6
1	C	730	LEU	2.6
1	B	734	SER	2.6
1	A	68	ALA	2.6
1	A	431	ARG	2.6
1	A	690	SER	2.6
1	D	734	SER	2.6
1	D	655	MET	2.6
1	A	76	CYS	2.5
1	D	131	GLU	2.5
1	B	686	PRO	2.4
1	B	735	HIS	2.4
1	D	74	LEU	2.4
1	A	136	GLU	2.4
1	D	130	ASP	2.4
1	D	1023	LYS	2.4
1	A	179	ALA	2.4
1	C	831	ALA	2.3
1	D	71	GLU	2.3
1	D	729	THR	2.3
1	D	179	ALA	2.3
1	D	128	ASN	2.3
1	C	663	LEU	2.3
1	D	633	GLY	2.3
1	A	129	VAL	2.3
1	C	689	GLU	2.3
1	B	688	PRO	2.3
1	B	729	THR	2.3
1	C	733	ALA	2.3
1	A	1023	LYS	2.3
1	A	72	SER	2.3
1	D	731	PRO	2.3
1	A	663	LEU	2.3
1	C	12	GLN	2.2
1	A	69	VAL	2.2
1	D	755	ARG	2.2
1	B	736	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	729	THR	2.2
1	D	79	PRO	2.2
1	C	772	ASP	2.2
1	A	180	GLY	2.2
1	D	690	SER	2.2
1	B	761	GLN	2.2
1	A	80	GLU	2.1
1	C	771	GLY	2.1
1	A	178	ARG	2.1
1	A	734	SER	2.1
1	C	734	SER	2.1
1	D	861	SER	2.1
1	D	745	MET	2.1
1	A	126	THR	2.1
1	D	699	ARG	2.1
1	D	78	LEU	2.1
1	C	581	ASN	2.1
1	A	117	GLU	2.1
1	A	181	GLU	2.1
1	B	745	MET	2.1
1	A	687	GLN	2.1
1	A	157	ARG	2.0
1	C	735	HIS	2.0
1	C	737	ILE	2.0
1	A	699	ARG	2.0
1	D	774	LYS	2.0
1	C	729	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	DMS	D	8009	4/4	0.66	0.38	78,79,80,80	0
4	DMS	A	8013	4/4	0.73	0.32	66,66,68,69	0
4	DMS	D	8014	4/4	0.74	0.29	69,69,71,73	0
4	DMS	D	8013	4/4	0.75	0.24	53,55,57,63	0
4	DMS	A	8009	4/4	0.75	0.31	67,70,71,72	0
4	DMS	C	8012	4/4	0.77	0.39	96,96,97,97	0
4	DMS	B	8012	4/4	0.78	0.34	63,64,65,66	0
4	DMS	D	8012	4/4	0.78	0.25	74,75,77,79	0
4	DMS	C	8017	4/4	0.79	0.32	63,63,63,64	0
4	DMS	A	8014	4/4	0.80	0.23	61,62,63,64	0
4	DMS	B	8011	4/4	0.80	0.25	65,67,68,69	0
4	DMS	A	8012	4/4	0.80	0.28	74,75,77,78	0
4	DMS	C	8016	4/4	0.82	0.24	58,60,60,62	0
4	DMS	A	8007	4/4	0.83	0.27	60,61,62,63	0
2	MG	A	3002	1/1	0.83	0.10	31,31,31,31	0
4	DMS	C	8013	4/4	0.83	0.28	58,60,61,62	0
4	DMS	A	8006	4/4	0.84	0.21	43,45,48,49	0
4	DMS	C	8009	4/4	0.84	0.22	47,50,51,56	0
4	DMS	B	8015	4/4	0.85	0.20	48,50,53,58	0
4	DMS	A	8017	4/4	0.86	0.18	54,56,56,61	0
4	DMS	B	8002	4/4	0.86	0.26	35,43,44,45	0
3	NA	C	3104	1/1	0.86	0.11	32,32,32,32	0
4	DMS	C	8011	4/4	0.86	0.23	53,53,53,54	0
4	DMS	C	8015	4/4	0.87	0.20	50,52,53,56	0
4	DMS	C	8003	4/4	0.87	0.20	46,50,54,54	0
4	DMS	C	8007	4/4	0.87	0.23	49,51,52,53	0
4	DMS	D	8002	4/4	0.87	0.26	42,48,50,52	0
4	DMS	C	8002	4/4	0.88	0.21	29,34,36,44	0
4	DMS	D	8004	4/4	0.88	0.16	36,37,42,46	0
4	DMS	D	8006	4/4	0.88	0.19	50,53,54,55	0
4	DMS	D	8007	4/4	0.88	0.25	59,59,60,61	0
4	DMS	B	8010	4/4	0.88	0.23	45,50,51,54	0
4	DMS	A	8002	4/4	0.88	0.23	40,40,42,48	0
4	DMS	A	8016	4/4	0.88	0.22	69,71,72,73	0
4	DMS	B	8007	4/4	0.88	0.22	45,47,50,53	0
4	DMS	B	8017	4/4	0.89	0.22	61,61,65,65	0
3	NA	B	3104	1/1	0.89	0.11	38,38,38,38	0
4	DMS	B	8008	4/4	0.89	0.19	42,44,44,48	0
4	DMS	C	8006	4/4	0.89	0.18	46,46,46,52	0
4	DMS	D	8010	4/4	0.89	0.20	46,49,51,51	0
4	DMS	B	8014	4/4	0.89	0.17	30,31,32,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	DMS	C	8008	4/4	0.89	0.21	53,54,55,56	0
4	DMS	B	8005	4/4	0.89	0.18	44,49,50,52	0
4	DMS	D	8015	4/4	0.89	0.18	50,51,53,55	0
4	DMS	A	8015	4/4	0.90	0.19	56,56,56,58	0
4	DMS	C	8019	4/4	0.90	0.18	60,61,61,65	0
4	DMS	A	8003	4/4	0.90	0.16	42,44,45,51	0
4	DMS	D	8001	4/4	0.91	0.18	35,38,38,46	0
4	DMS	B	8004	4/4	0.91	0.17	34,37,38,42	0
4	DMS	C	8004	4/4	0.91	0.17	32,36,37,39	0
4	DMS	D	8005	4/4	0.91	0.19	35,39,41,43	0
4	DMS	B	8013	4/4	0.91	0.20	56,56,57,58	0
4	DMS	C	8014	4/4	0.91	0.19	65,67,67,70	0
2	MG	D	3002	1/1	0.92	0.09	28,28,28,28	0
3	NA	B	3103	1/1	0.92	0.12	37,37,37,37	0
4	DMS	B	8003	4/4	0.92	0.13	29,30,37,41	0
2	MG	B	3002	1/1	0.92	0.10	17,17,17,17	0
4	DMS	A	8004	4/4	0.92	0.15	43,47,49,50	0
4	DMS	B	8006	4/4	0.92	0.12	36,36,38,43	0
3	NA	B	3102	1/1	0.93	0.07	20,20,20,20	0
4	DMS	B	8001	4/4	0.93	0.16	31,32,35,38	0
4	DMS	C	8001	4/4	0.93	0.15	27,31,34,37	0
4	DMS	C	8005	4/4	0.93	0.18	51,51,53,55	0
5	IPT	D	2001	15/15	0.93	0.10	26,29,33,34	0
4	DMS	A	8008	4/4	0.95	0.12	43,43,46,47	0
4	DMS	C	8010	4/4	0.95	0.12	40,40,43,45	0
4	DMS	D	8003	4/4	0.95	0.11	37,38,38,39	0
3	NA	A	3103	1/1	0.95	0.06	36,36,36,36	0
4	DMS	A	8010	4/4	0.95	0.13	42,42,42,43	0
4	DMS	C	8018	4/4	0.95	0.11	29,34,36,37	0
3	NA	A	3102	1/1	0.95	0.04	17,17,17,17	0
4	DMS	A	8011	4/4	0.96	0.12	40,44,45,47	0
4	DMS	B	8016	4/4	0.96	0.10	40,40,42,42	0
2	MG	D	3001	1/1	0.96	0.04	26,26,26,26	0
2	MG	C	3002	1/1	0.96	0.09	15,15,15,15	0
3	NA	A	3101	1/1	0.96	0.08	26,26,26,26	0
5	IPT	A	2001	15/15	0.96	0.08	20,23,32,32	0
5	IPT	B	2001	15/15	0.96	0.07	15,21,26,28	0
5	IPT	C	2001	15/15	0.96	0.07	16,21,32,35	0
3	NA	D	3103	1/1	0.96	0.06	30,30,30,30	0
4	DMS	D	8011	4/4	0.97	0.11	37,40,42,42	0
2	MG	B	3007	1/1	0.97	0.04	35,35,35,35	0
3	NA	D	3101	1/1	0.97	0.06	24,24,24,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	DMS	A	8005	4/4	0.97	0.09	29,30,30,36	0
3	NA	D	3102	1/1	0.97	0.06	18,18,18,18	0
2	MG	A	3001	1/1	0.97	0.04	25,25,25,25	0
4	DMS	D	8008	4/4	0.97	0.10	48,48,48,50	0
4	DMS	B	8009	4/4	0.97	0.10	32,37,38,41	0
3	NA	C	3102	1/1	0.97	0.04	18,18,18,18	0
4	DMS	A	8001	4/4	0.98	0.09	27,33,33,34	0
3	NA	B	3101	1/1	0.98	0.04	16,16,16,16	0
3	NA	C	3103	1/1	0.98	0.05	35,35,35,35	0
2	MG	B	3001	1/1	0.98	0.03	21,21,21,21	0
2	MG	C	3001	1/1	0.99	0.04	19,19,19,19	0
3	NA	C	3101	1/1	0.99	0.02	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.