



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 7, 2020 – 07:07 PM BST

PDB ID : 1JZ4  
Title : E. COLI (lacZ) BETA-GALACTOSIDASE-TRAPPED 2-DEOXY-GALACTOSYL-ENZYME INTERMEDIATE (Low Bis-Tris)  
Authors : Juers, D.H.; Matthews, B.W.  
Deposited on : 2001-09-13  
Resolution : 2.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

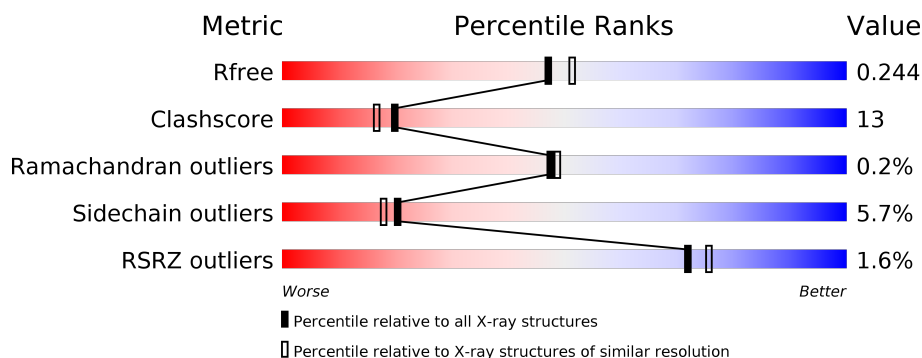
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	1023	<div> <div>2%</div> <div> <div></div> <div>64%</div> <div>29%</div> <div>5% ..</div> </div> </div>
1	B	1023	<div> <div>%</div> <div> <div></div> <div>62%</div> <div>30%</div> <div>5% ..</div> </div> </div>
1	C	1023	<div> <div>2%</div> <div> <div></div> <div>63%</div> <div>28%</div> <div>7% ..</div> </div> </div>
1	D	1023	<div> <div>2%</div> <div> <div></div> <div>63%</div> <div>30%</div> <div>5% ..</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	DMS	B	8508	-	X	-	-
5	DMS	C	8506	-	-	X	-
5	DMS	D	8412	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 36890 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Beta-Galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1011	Total	C	N	O	S	0	0	0
			8125	5138	1440	1509	38			
1	B	1011	Total	C	N	O	S	0	0	0
			8125	5138	1440	1509	38			
1	C	1011	Total	C	N	O	S	0	0	0
			8125	5138	1440	1509	38			
1	D	1011	Total	C	N	O	S	0	0	0
			8125	5138	1440	1509	38			

There are 32 discrepancies between the modelled and reference sequences:

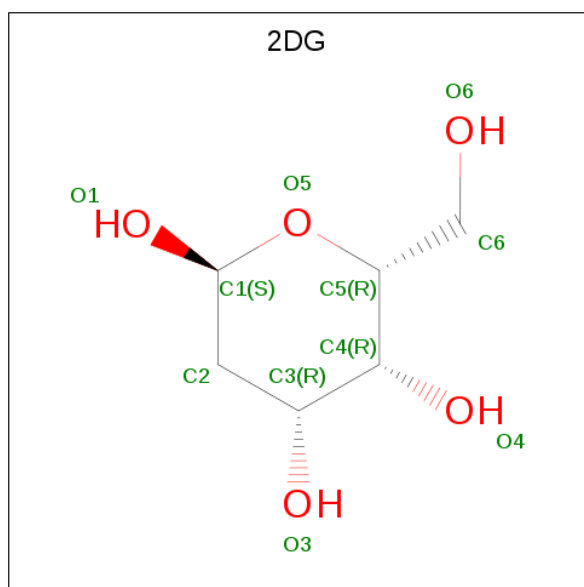
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	THR	cloning artifact	? P00722
A	2	SER	MET	cloning artifact	? P00722
A	3	HIS	ILE	cloning artifact	? P00722
A	4	MET	THR	cloning artifact	? P00722
A	5	LEU	ASP	cloning artifact	? P00722
A	6	GLU	SER	cloning artifact	? P00722
A	7	ASP	LEU	cloning artifact	? P00722
A	8	PRO	ALA	cloning artifact	? P00722
B	1	GLY	THR	cloning artifact	? P00722
B	2	SER	MET	cloning artifact	? P00722
B	3	HIS	ILE	cloning artifact	? P00722
B	4	MET	THR	cloning artifact	? P00722
B	5	LEU	ASP	cloning artifact	? P00722
B	6	GLU	SER	cloning artifact	? P00722
B	7	ASP	LEU	cloning artifact	? P00722
B	8	PRO	ALA	cloning artifact	? P00722
C	1	GLY	THR	cloning artifact	? P00722
C	2	SER	MET	cloning artifact	? P00722
C	3	HIS	ILE	cloning artifact	? P00722
C	4	MET	THR	cloning artifact	? P00722
C	5	LEU	ASP	cloning artifact	? P00722

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Chain	Residue	Modelled	Actual	Comment	Reference
C	6	GLU	SER	cloning artifact	? P00722
C	7	ASP	LEU	cloning artifact	? P00722
C	8	PRO	ALA	cloning artifact	? P00722
D	1	GLY	THR	cloning artifact	? P00722
D	2	SER	MET	cloning artifact	? P00722
D	3	HIS	ILE	cloning artifact	? P00722
D	4	MET	THR	cloning artifact	? P00722
D	5	LEU	ASP	cloning artifact	? P00722
D	6	GLU	SER	cloning artifact	? P00722
D	7	ASP	LEU	cloning artifact	? P00722
D	8	PRO	ALA	cloning artifact	? P00722

- Molecule 2 is 2-deoxy-alpha-D-galactopyranose (three-letter code: 2DG) (formula:  $C_6H_{12}O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		
2	B	1	Total	C	O	0	0
			10	6	4		
2	C	1	Total	C	O	0	0
			10	6	4		
2	D	1	Total	C	O	0	0
			10	6	4		

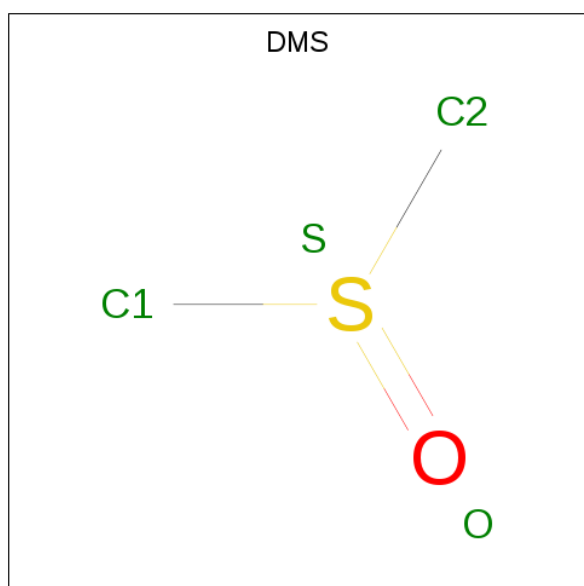
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C<sub>2</sub>H<sub>6</sub>OS).



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

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

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

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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
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.

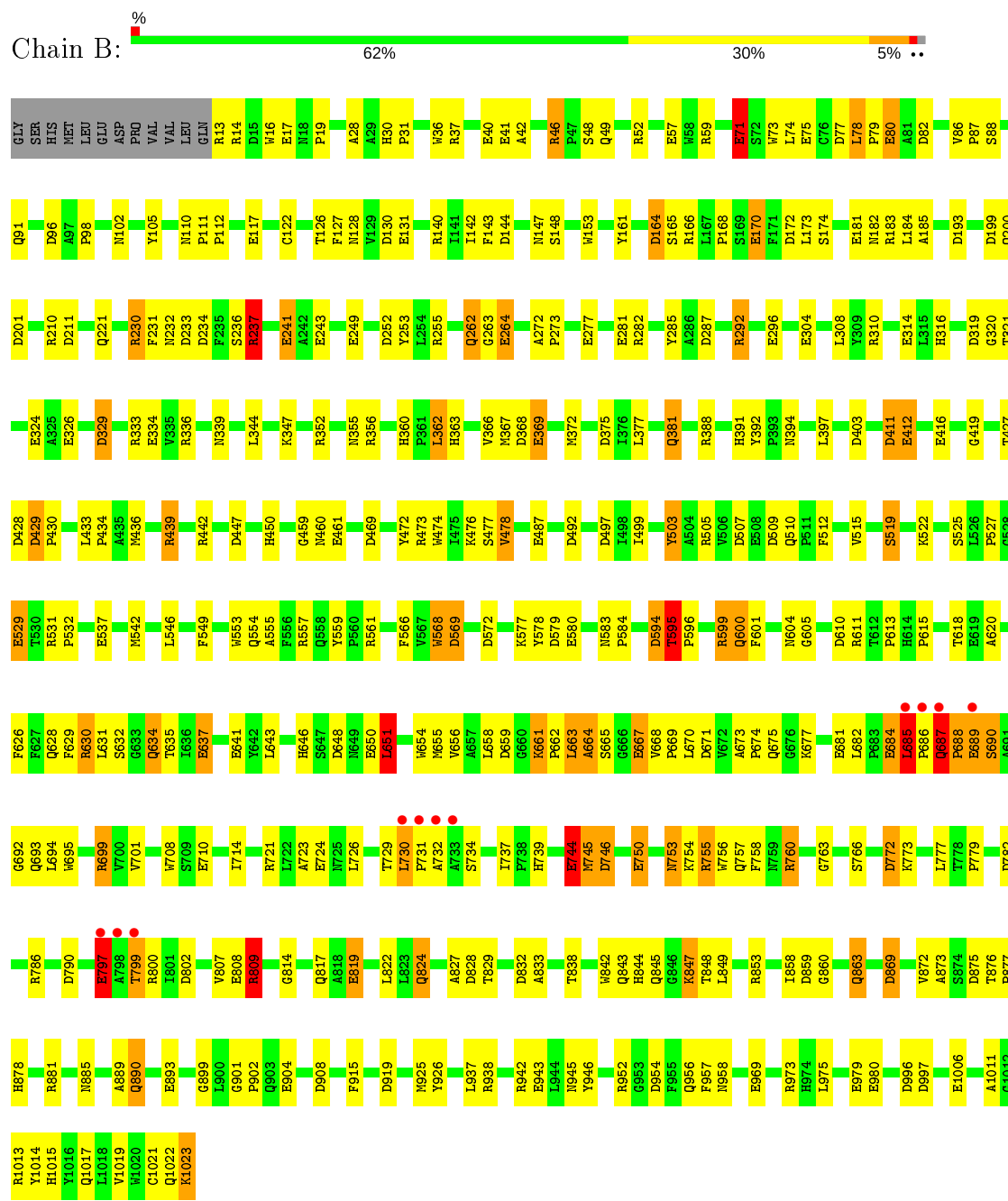
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	992	Total	O	0	0
			992	992		
6	B	995	Total	O	0	0
			995	995		

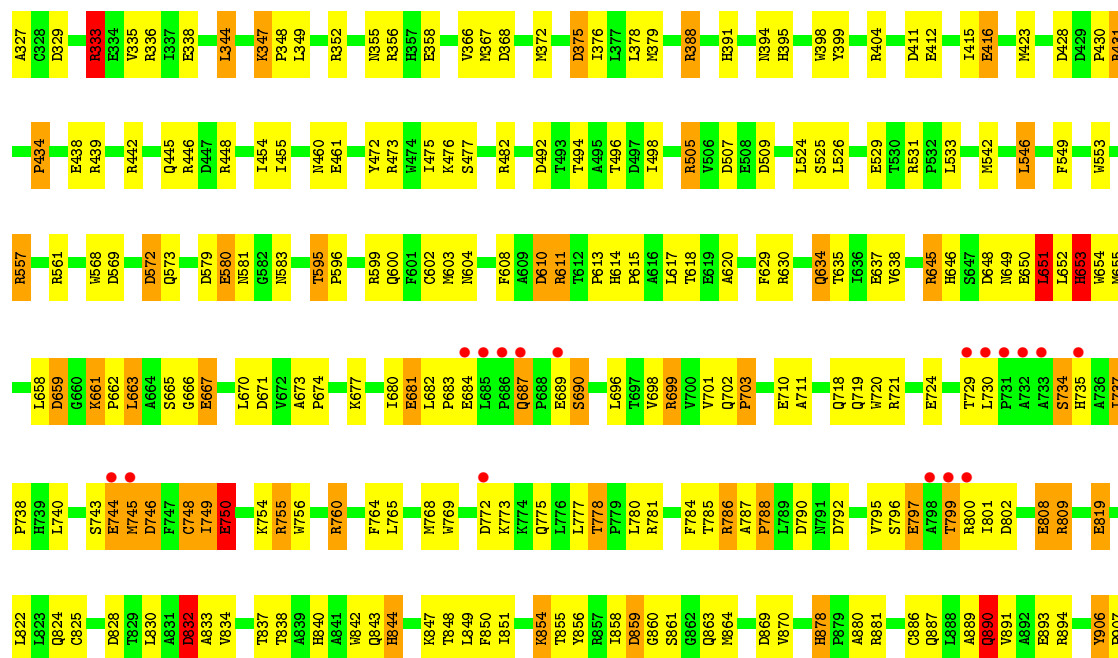
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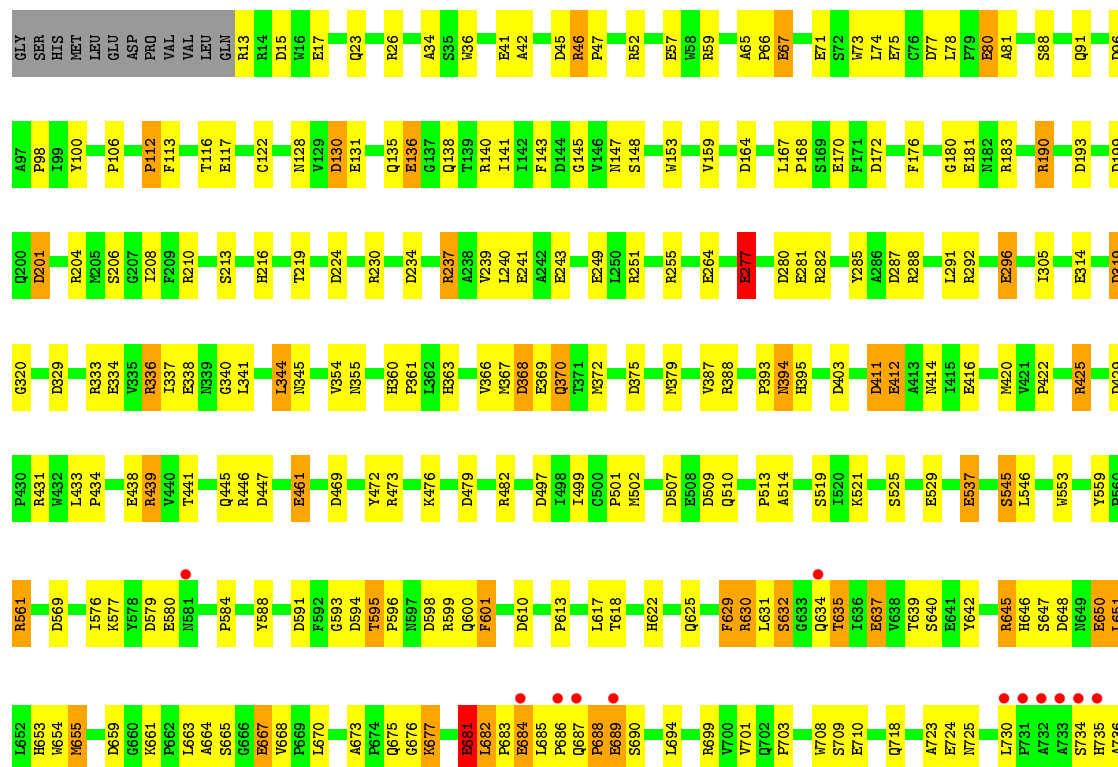
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	946	Total 946	O 946	0	0
6	D	980	Total 980	O 980	0	0

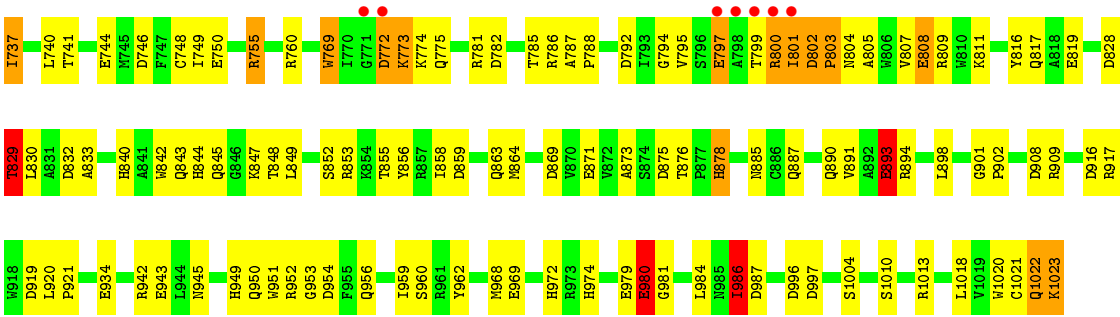






### • Molecule 1: Beta-Galactosidase







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	149.50Å 169.00Å 200.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.10 19.99 – 2.10	Depositor EDS
% Data completeness (in resolution range)	90.0 (40.00-2.10) 86.4 (19.99-2.10)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.18 (at 2.09Å)	Xtriage
Refinement program	TNT	Depositor
R, $R_{free}$	0.164 , 0.267 0.152 , 0.244	Depositor DCC
$R_{free}$ test set	3827 reflections (1.44%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	11.5	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 95.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	36890	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	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 36.06 % 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 5.3009e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NA, MG, DMS, 2DG

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	1.09	43/8367 (0.5%)	1.73	174/11415 (1.5%)
1	B	1.11	48/8367 (0.6%)	1.69	165/11415 (1.4%)
1	C	1.11	42/8367 (0.5%)	1.75	177/11415 (1.6%)
1	D	1.11	48/8367 (0.6%)	1.69	151/11415 (1.3%)
All	All	1.10	181/33468 (0.5%)	1.71	667/45660 (1.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	D	1	0
All	All	1	1

The worst 5 of 181 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	277	GLU	CD-OE2	10.41	1.37	1.25
1	C	281	GLU	CD-OE2	10.15	1.36	1.25
1	D	71	GLU	CD-OE2	10.14	1.36	1.25
1	C	684	GLU	CD-OE2	9.88	1.36	1.25
1	C	296	GLU	CD-OE2	9.86	1.36	1.25

The worst 5 of 667 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	809	ARG	NE-CZ-NH1	19.72	130.16	120.30
1	C	442	ARG	NE-CZ-NH2	-17.27	111.67	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	786	ARG	NE-CZ-NH1	15.02	127.81	120.30
1	C	721	ARG	NE-CZ-NH1	15.00	127.80	120.30
1	A	46	ARG	NE-CZ-NH1	15.00	127.80	120.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	D	689	GLU	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	329	ASP	Sidechain

## 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	8125	0	7716	188	1
1	B	8125	0	7716	201	0
1	C	8125	0	7716	219	0
1	D	8125	0	7716	208	0
2	A	10	0	9	0	0
2	B	10	0	9	1	0
2	C	10	0	9	0	0
2	D	10	0	9	2	0
3	A	3	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	3	0	0	0	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
4	C	4	0	0	0	0
4	D	3	0	0	0	0
5	A	100	0	150	9	0
5	B	108	0	162	10	0
5	C	112	0	168	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	92	0	138	13	0
6	A	992	0	0	21	0
6	B	995	0	0	14	0
6	C	946	0	0	14	1
6	D	980	0	0	18	0
All	All	36890	0	31518	810	1

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 13.

The worst 5 of 810 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:C:634:GLN:NE2	1:C:634:GLN:H	1.47	1.13
1:C:746:ASP:HA	1:C:760:ARG:HG3	1.33	1.10
1:D:804:ASN:HD22	1:D:809:ARG:NH2	1.54	1.05
1:B:142:ILE:HG12	1:B:170:GLU:HG2	1.37	1.02
1:B:797:GLU:HG2	1:B:799:THR:HG23	1.35	1.02

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:773:LYS:NZ	6:C:9410:HOH:O[3_544]	2.18	0.02

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	1009/1023 (99%)	964 (96%)	44 (4%)	1 (0%)	51 54
1	B	1009/1023 (99%)	956 (95%)	48 (5%)	5 (0%)	29 26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	1009/1023 (99%)	963 (95%)	46 (5%)	0	100	100
1	D	1009/1023 (99%)	962 (95%)	43 (4%)	4 (0%)	34	32
All	All	4036/4092 (99%)	3845 (95%)	181 (4%)	10 (0%)	47	49

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	690	SER
1	B	731	PRO
1	D	688	PRO
1	D	801	ILE
1	B	688	PRO

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	864/875 (99%)	816 (94%)	48 (6%)	21	18
1	B	864/875 (99%)	815 (94%)	49 (6%)	20	18
1	C	864/875 (99%)	817 (95%)	47 (5%)	22	20
1	D	864/875 (99%)	812 (94%)	52 (6%)	19	16
All	All	3456/3500 (99%)	3260 (94%)	196 (6%)	20	18

5 of 196 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	824	GLN
1	C	333	ARG
1	D	772	ASP
1	B	847	LYS
1	C	75	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 50 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	646	HIS
1	C	102	ASN
1	D	878	HIS
1	B	757	GLN
1	B	824	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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 132 ligands modelled in this entry, 25 are monoatomic - leaving 107 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	8412	-	3,3,3	1.92	1 (33%)	3,3,3	0.18	0
5	DMS	B	8423	-	3,3,3	0.57	0	3,3,3	0.15	0
5	DMS	A	8412	-	3,3,3	0.76	0	3,3,3	0.20	0
5	DMS	C	8602	-	3,3,3	1.06	0	3,3,3	0.16	0
5	DMS	A	8403	-	3,3,3	0.97	0	3,3,3	0.74	0
5	DMS	B	8425	4	3,3,3	1.92	1 (33%)	3,3,3	0.37	0
5	DMS	B	8417	-	3,3,3	1.18	0	3,3,3	0.48	0
5	DMS	B	8403	-	3,3,3	1.46	0	3,3,3	1.25	1 (33%)
5	DMS	D	8402	-	3,3,3	1.15	0	3,3,3	0.63	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	DMS	C	8408	-	3,3,3	0.68	0	3,3,3	0.25	0
5	DMS	A	8405	-	3,3,3	1.50	1 (33%)	3,3,3	0.24	0
5	DMS	A	8409	-	3,3,3	2.34	2 (66%)	3,3,3	0.36	0
2	2DG	B	2001	1,4	10,10,11	1.41	1 (10%)	13,13,15	1.27	1 (7%)
5	DMS	A	8406	-	3,3,3	1.88	1 (33%)	3,3,3	0.45	0
5	DMS	A	8413	-	3,3,3	2.03	2 (66%)	3,3,3	0.16	0
5	DMS	C	8427	-	3,3,3	0.90	0	3,3,3	0.09	0
5	DMS	A	8402	-	3,3,3	1.57	1 (33%)	3,3,3	0.14	0
5	DMS	B	8410	-	3,3,3	1.30	0	3,3,3	0.49	0
5	DMS	C	8503	-	3,3,3	0.79	0	3,3,3	0.42	0
5	DMS	B	8421	-	3,3,3	0.84	0	3,3,3	0.14	0
5	DMS	D	8421	-	3,3,3	1.03	0	3,3,3	0.08	0
5	DMS	B	8415	-	3,3,3	1.10	0	3,3,3	1.33	1 (33%)
5	DMS	D	8401	-	3,3,3	1.78	1 (33%)	3,3,3	0.33	0
5	DMS	D	8416	-	3,3,3	0.77	0	3,3,3	0.23	0
5	DMS	C	8415	-	3,3,3	1.57	1 (33%)	3,3,3	0.97	0
5	DMS	D	8413	-	3,3,3	0.89	0	3,3,3	0.65	0
5	DMS	C	8420	-	3,3,3	1.29	0	3,3,3	0.66	0
5	DMS	D	8503	-	3,3,3	1.34	1 (33%)	3,3,3	0.78	0
5	DMS	A	8602	-	3,3,3	0.87	0	3,3,3	0.83	0
5	DMS	C	8411	-	3,3,3	1.68	1 (33%)	3,3,3	0.34	0
5	DMS	A	8502	-	3,3,3	1.16	0	3,3,3	1.36	1 (33%)
5	DMS	D	8501	-	3,3,3	0.84	0	3,3,3	0.90	0
5	DMS	D	8406	-	3,3,3	1.02	0	3,3,3	0.38	0
5	DMS	B	8405	-	3,3,3	0.89	0	3,3,3	0.39	0
5	DMS	A	8503	-	3,3,3	0.66	0	3,3,3	0.18	0
5	DMS	D	8405	-	3,3,3	0.83	0	3,3,3	0.25	0
5	DMS	A	8401	-	3,3,3	1.94	1 (33%)	3,3,3	0.12	0
5	DMS	C	8425	4	3,3,3	1.48	1 (33%)	3,3,3	1.65	1 (33%)
5	DMS	C	8504	-	3,3,3	1.11	0	3,3,3	0.23	0
5	DMS	C	8410	-	3,3,3	1.10	0	3,3,3	0.36	0
5	DMS	C	8413	-	3,3,3	2.95	2 (66%)	3,3,3	0.43	0
5	DMS	C	8601	-	3,3,3	1.09	0	3,3,3	0.61	0
5	DMS	B	8408	-	3,3,3	1.41	0	3,3,3	0.40	0
5	DMS	B	8404	-	3,3,3	0.81	0	3,3,3	0.36	0
5	DMS	A	8404	-	3,3,3	1.82	1 (33%)	3,3,3	0.31	0
2	2DG	A	2001	1,4	10,10,11	1.04	1 (10%)	13,13,15	1.89	3 (23%)
5	DMS	C	8501	-	3,3,3	1.53	0	3,3,3	1.19	1 (33%)
5	DMS	D	8411	-	3,3,3	0.93	0	3,3,3	0.49	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	2DG	C	2001	1,4	10,10,11	1.28	1 (10%)	13,13,15	1.87	1 (7%)
5	DMS	B	8416	-	3,3,3	1.30	0	3,3,3	0.07	0
5	DMS	B	8504	-	3,3,3	0.11	0	3,3,3	0.08	0
5	DMS	B	8427	-	3,3,3	0.88	0	3,3,3	0.61	0
5	DMS	D	8403	-	3,3,3	0.52	0	3,3,3	0.31	0
5	DMS	B	8502	-	3,3,3	0.74	0	3,3,3	1.26	1 (33%)
5	DMS	C	8407	-	3,3,3	1.16	0	3,3,3	0.11	0
5	DMS	B	8409	-	3,3,3	0.71	0	3,3,3	1.28	1 (33%)
5	DMS	D	8408	-	3,3,3	1.14	0	3,3,3	0.32	0
5	DMS	C	8409	-	3,3,3	1.58	1 (33%)	3,3,3	0.38	0
5	DMS	B	8407	-	3,3,3	1.51	0	3,3,3	0.26	0
5	DMS	A	8414	-	3,3,3	0.76	0	3,3,3	0.44	0
5	DMS	A	8504	-	3,3,3	0.35	0	3,3,3	0.42	0
5	DMS	D	8412	-	3,3,3	1.49	1 (33%)	3,3,3	0.59	0
5	DMS	C	8405	-	3,3,3	1.83	1 (33%)	3,3,3	0.26	0
5	DMS	B	8420	-	3,3,3	1.30	1 (33%)	3,3,3	0.39	0
5	DMS	A	8415	-	3,3,3	1.94	1 (33%)	3,3,3	0.15	0
5	DMS	A	8411	-	3,3,3	1.56	1 (33%)	3,3,3	0.15	0
5	DMS	A	8410	-	3,3,3	0.72	0	3,3,3	0.09	0
5	DMS	A	8501	-	3,3,3	1.45	1 (33%)	3,3,3	0.32	0
5	DMS	C	8404	-	3,3,3	1.29	0	3,3,3	0.68	0
5	DMS	B	8601	-	3,3,3	0.43	0	3,3,3	0.29	0
5	DMS	C	8401	-	3,3,3	0.88	0	3,3,3	0.76	0
5	DMS	C	8402	-	3,3,3	2.08	1 (33%)	3,3,3	0.72	0
5	DMS	B	8414	-	3,3,3	0.56	0	3,3,3	0.76	0
5	DMS	D	8701	-	3,3,3	2.89	2 (66%)	3,3,3	0.76	0
5	DMS	D	8417	-	3,3,3	1.27	0	3,3,3	0.72	0
5	DMS	B	8508	-	3,3,3	2.50	3 (100%)	3,3,3	0.38	0
5	DMS	B	8506	-	3,3,3	1.39	0	3,3,3	0.79	0
5	DMS	C	8416	-	3,3,3	1.11	0	3,3,3	0.72	0
5	DMS	D	8508	-	3,3,3	0.56	0	3,3,3	0.59	0
5	DMS	D	8419	-	3,3,3	0.88	0	3,3,3	0.36	0
5	DMS	D	8705	-	3,3,3	1.23	0	3,3,3	0.21	0
5	DMS	D	8414	-	3,3,3	1.22	0	3,3,3	0.08	0
5	DMS	C	8414	-	3,3,3	1.59	1 (33%)	3,3,3	0.36	0
5	DMS	B	8406	-	3,3,3	1.26	1 (33%)	3,3,3	0.26	0
5	DMS	C	8417	-	3,3,3	0.86	0	3,3,3	0.52	0
5	DMS	C	8403	-	3,3,3	0.84	0	3,3,3	0.43	0
5	DMS	B	8402	-	3,3,3	1.89	2 (66%)	3,3,3	0.18	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	DMS	C	8412	-	3,3,3	0.87	0	3,3,3	0.56	0
5	DMS	A	8425	4	3,3,3	1.42	1 (33%)	3,3,3	0.56	0
5	DMS	D	8404	-	3,3,3	1.58	1 (33%)	3,3,3	0.51	0
5	DMS	A	8419	-	3,3,3	0.66	0	3,3,3	0.40	0
5	DMS	A	8417	-	3,3,3	1.12	0	3,3,3	0.50	0
5	DMS	C	8506	-	3,3,3	1.59	1 (33%)	3,3,3	1.25	1 (33%)
5	DMS	A	8420	-	3,3,3	1.57	1 (33%)	3,3,3	1.02	0
5	DMS	B	8401	-	3,3,3	0.66	0	3,3,3	0.44	0
5	DMS	D	8703	-	3,3,3	0.84	0	3,3,3	0.33	0
5	DMS	D	8410	-	3,3,3	1.30	0	3,3,3	0.49	0
2	2DG	D	2001	1,4	10,10,11	1.42	2 (20%)	13,13,15	2.35	4 (30%)
5	DMS	C	8419	-	3,3,3	1.40	0	3,3,3	0.19	0
5	DMS	D	8409	-	3,3,3	2.22	1 (33%)	3,3,3	0.59	0
5	DMS	C	8423	-	3,3,3	1.30	0	3,3,3	0.31	0
5	DMS	B	8411	-	3,3,3	0.36	0	3,3,3	0.76	0
5	DMS	A	8408	-	3,3,3	1.06	0	3,3,3	0.24	0
5	DMS	A	8421	-	3,3,3	0.89	0	3,3,3	0.97	0
5	DMS	B	8413	-	3,3,3	2.23	2 (66%)	3,3,3	0.82	0
5	DMS	C	8421	-	3,3,3	0.60	0	3,3,3	0.23	0
5	DMS	A	8407	-	3,3,3	1.74	2 (66%)	3,3,3	0.54	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	2DG	C	2001	1,4	-	1/2/16/18	0/1/1/1
2	2DG	D	2001	1,4	-	1/2/16/18	0/1/1/1
2	2DG	A	2001	1,4	-	1/2/16/18	0/1/1/1
2	2DG	B	2001	1,4	-	1/2/16/18	0/1/1/1

The worst 5 of 49 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	8413	DMS	O-S	4.50	1.80	1.50
5	D	8701	DMS	O-S	3.75	1.75	1.50
5	D	8409	DMS	O-S	3.74	1.75	1.50
2	D	2001	2DG	C2-C3	3.40	1.57	1.52
5	A	8409	DMS	O-S	3.29	1.72	1.50

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2001	2DG	C3-C4-C5	-5.77	104.22	109.97
2	A	2001	2DG	C3-C4-C5	-4.74	105.25	109.97
2	D	2001	2DG	C1-O5-C5	4.27	119.43	112.01
2	D	2001	2DG	C2-C3-C4	4.25	116.04	111.16
2	D	2001	2DG	C3-C4-C5	-4.21	105.77	109.97

There are no chirality outliers.

All (4) torsion outliers are listed below:

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

There are no ring outliers.

31 monomers are involved in 50 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	8412	DMS	1	0
5	C	8602	DMS	1	0
5	B	8417	DMS	1	0
2	B	2001	2DG	1	0
5	A	8406	DMS	1	0
5	C	8427	DMS	1	0
5	C	8503	DMS	3	0
5	B	8415	DMS	2	0
5	D	8416	DMS	3	0
5	D	8413	DMS	1	0
5	C	8420	DMS	2	0
5	D	8503	DMS	2	0
5	A	8602	DMS	1	0
5	A	8502	DMS	3	0
5	C	8504	DMS	1	0
5	C	8601	DMS	1	0
5	B	8504	DMS	1	0
5	A	8414	DMS	1	0
5	D	8412	DMS	4	0
5	A	8410	DMS	1	0
5	B	8601	DMS	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	8506	DMS	2	0
5	D	8508	DMS	1	0
5	D	8419	DMS	1	0
5	B	8406	DMS	1	0
5	B	8402	DMS	1	0
5	C	8412	DMS	2	0
5	C	8506	DMS	4	0
5	A	8420	DMS	2	0
5	D	8703	DMS	1	0
2	D	2001	2DG	2	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	1011/1023 (98%)	-0.76	16 (1%) 72 75	5, 15, 44, 98	0
1	B	1011/1023 (98%)	-0.73	11 (1%) 80 84	5, 15, 45, 100	0
1	C	1011/1023 (98%)	-0.70	17 (1%) 70 74	6, 15, 46, 100	0
1	D	1011/1023 (98%)	-0.70	19 (1%) 66 71	4, 16, 46, 99	0
All	All	4044/4092 (98%)	-0.72	63 (1%) 72 75	4, 15, 45, 100	0

The worst 5 of 63 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	731	PRO	8.2
1	A	735	HIS	8.0
1	A	686	PRO	7.8
1	B	731	PRO	7.3
1	D	732	ALA	6.5

### 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 monosaccharides 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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	DMS	D	8703	4/4	0.89	0.21	24,53,80,80	0
5	DMS	B	8508	4/4	0.90	0.11	27,31,46,64	0
3	MG	A	3005	1/1	0.90	0.04	36,36,36,36	0
4	NA	D	3103	1/1	0.91	0.11	38,38,38,38	0
5	DMS	B	8406	4/4	0.92	0.19	34,40,61,78	0
5	DMS	A	8502	4/4	0.93	0.12	14,21,75,84	0
5	DMS	A	8417	4/4	0.93	0.17	25,26,61,100	0
5	DMS	B	8427	4/4	0.93	0.10	28,31,42,65	0
5	DMS	B	8413	4/4	0.93	0.15	22,32,35,42	0
5	DMS	A	8415	4/4	0.94	0.10	24,27,28,70	0
5	DMS	D	8417	4/4	0.94	0.14	18,26,39,73	0
5	DMS	B	8408	4/4	0.94	0.17	15,25,43,60	0
5	DMS	C	8419	4/4	0.94	0.14	33,44,54,61	0
5	DMS	B	8415	4/4	0.94	0.12	22,26,33,67	0
5	DMS	B	8407	4/4	0.95	0.17	34,34,63,66	0
4	NA	B	3104	1/1	0.95	0.09	37,37,37,37	0
5	DMS	C	8417	4/4	0.95	0.10	24,51,57,58	0
5	DMS	A	8425	4/4	0.95	0.15	34,40,45,45	0
5	DMS	C	8413	4/4	0.95	0.16	31,32,34,38	0
5	DMS	C	8415	4/4	0.95	0.11	20,26,33,52	0
5	DMS	B	8425	4/4	0.95	0.12	22,30,32,54	0
5	DMS	C	8423	4/4	0.95	0.10	26,27,32,61	0
5	DMS	A	8421	4/4	0.95	0.18	42,43,67,100	0
5	DMS	C	8602	4/4	0.95	0.11	27,32,53,77	0
5	DMS	C	8425	4/4	0.96	0.11	33,37,39,48	0
5	DMS	A	8409	4/4	0.96	0.11	24,27,39,42	0
5	DMS	C	8601	4/4	0.96	0.14	31,53,55,100	0
5	DMS	C	8416	4/4	0.96	0.26	31,33,49,64	0
5	DMS	D	8705	4/4	0.96	0.12	24,32,51,51	0
5	DMS	C	8503	4/4	0.96	0.19	20,31,54,100	0
5	DMS	A	8406	4/4	0.96	0.13	7,53,55,100	0
5	DMS	B	8421	4/4	0.96	0.10	14,46,60,85	0
5	DMS	B	8502	4/4	0.96	0.09	13,30,45,50	0
5	DMS	C	8506	4/4	0.96	0.10	15,24,47,89	0
5	DMS	B	8412	4/4	0.96	0.11	21,33,34,38	0
2	2DG	D	2001	10/11	0.96	0.08	9,15,24,24	0
5	DMS	B	8423	4/4	0.96	0.10	28,44,71,87	0
5	DMS	D	8409	4/4	0.96	0.07	25,30,39,53	0
5	DMS	B	8420	4/4	0.96	0.13	36,45,63,100	0
5	DMS	B	8417	4/4	0.96	0.10	26,31,34,63	0
5	DMS	D	8406	4/4	0.96	0.08	18,19,20,34	0
5	DMS	A	8407	4/4	0.96	0.13	28,40,42,43	0
2	2DG	A	2001	10/11	0.97	0.07	10,15,18,25	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	2DG	C	2001	10/11	0.97	0.08	7,15,26,27	0
5	DMS	B	8416	4/4	0.97	0.12	35,42,52,88	0
4	NA	C	3103	1/1	0.97	0.06	22,22,22,22	0
2	2DG	B	2001	10/11	0.97	0.11	13,15,19,24	0
5	DMS	D	8414	4/4	0.97	0.14	17,32,44,92	0
4	NA	C	3104	1/1	0.97	0.08	26,26,26,26	0
5	DMS	D	8501	4/4	0.97	0.08	16,34,36,46	0
5	DMS	B	8409	4/4	0.97	0.09	18,32,33,38	0
5	DMS	C	8409	4/4	0.97	0.09	23,35,41,41	0
5	DMS	D	8421	4/4	0.97	0.11	36,45,48,52	0
5	DMS	A	8503	4/4	0.97	0.20	42,47,48,100	0
5	DMS	A	8413	4/4	0.97	0.12	34,34,53,100	0
5	DMS	D	8416	4/4	0.97	0.21	16,34,42,89	0
3	MG	D	3005	1/1	0.97	0.05	43,43,43,43	0
5	DMS	A	8501	4/4	0.97	0.08	12,13,29,33	0
5	DMS	B	8601	4/4	0.97	0.11	36,60,100,100	0
5	DMS	B	8414	4/4	0.97	0.13	24,43,49,100	0
5	DMS	D	8503	4/4	0.97	0.18	33,53,64,100	0
5	DMS	A	8412	4/4	0.98	0.10	24,33,37,37	0
3	MG	B	3001	1/1	0.98	0.06	11,11,11,11	0
5	DMS	D	8701	4/4	0.98	0.09	15,17,22,37	0
5	DMS	B	8504	4/4	0.98	0.10	31,37,51,52	0
5	DMS	C	8501	4/4	0.98	0.12	23,23,34,38	0
5	DMS	B	8506	4/4	0.98	0.10	31,39,63,100	0
4	NA	C	3101	1/1	0.98	0.05	11,11,11,11	0
5	DMS	D	8413	4/4	0.98	0.13	21,24,34,35	0
5	DMS	C	8420	4/4	0.98	0.12	34,36,53,100	0
4	NA	A	3103	1/1	0.98	0.11	25,25,25,25	0
5	DMS	D	8508	4/4	0.98	0.08	19,46,50,80	0
5	DMS	D	8419	4/4	0.98	0.07	18,32,40,44	0
5	DMS	A	8602	4/4	0.98	0.18	42,45,100,100	0
5	DMS	C	8427	4/4	0.98	0.15	37,43,81,82	0
5	DMS	C	8414	4/4	0.98	0.08	17,18,23,39	0
5	DMS	C	8407	4/4	0.98	0.11	35,39,51,52	0
5	DMS	A	8402	4/4	0.98	0.07	11,26,30,47	0
5	DMS	B	8402	4/4	0.98	0.07	8,21,25,29	0
5	DMS	B	8410	4/4	0.98	0.09	18,27,41,54	0
5	DMS	A	8419	4/4	0.98	0.12	24,26,55,100	0
5	DMS	D	8402	4/4	0.98	0.07	9,23,27,32	0
5	DMS	A	8414	4/4	0.98	0.10	16,37,47,100	0
4	NA	A	3101	1/1	0.98	0.05	13,13,13,13	0
5	DMS	A	8420	4/4	0.98	0.08	33,35,39,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	DMS	C	8408	4/4	0.98	0.09	13,35,46,83	0
4	NA	B	3101	1/1	0.98	0.06	13,13,13,13	0
5	DMS	D	8410	4/4	0.98	0.10	22,35,39,100	0
5	DMS	A	8504	4/4	0.98	0.10	23,35,38,100	0
5	DMS	D	8412	4/4	0.98	0.10	15,25,32,33	0
4	NA	B	3102	1/1	0.98	0.06	13,13,13,13	0
3	MG	D	3002	1/1	0.98	0.07	19,19,19,19	0
5	DMS	A	8411	4/4	0.98	0.08	24,26,28,100	0
5	DMS	C	8504	4/4	0.98	0.07	32,34,42,64	0
5	DMS	C	8421	4/4	0.98	0.11	26,26,41,64	0
3	MG	A	3002	1/1	0.98	0.05	14,14,14,14	0
4	NA	D	3102	1/1	0.99	0.05	18,18,18,18	0
4	NA	A	3102	1/1	0.99	0.04	15,15,15,15	0
5	DMS	D	8408	4/4	0.99	0.10	14,27,28,36	0
5	DMS	B	8405	4/4	0.99	0.11	23,28,62,71	0
4	NA	B	3103	1/1	0.99	0.03	19,19,19,19	0
5	DMS	C	8410	4/4	0.99	0.07	21,35,37,43	0
5	DMS	C	8411	4/4	0.99	0.08	18,20,34,50	0
3	MG	C	3002	1/1	0.99	0.06	12,12,12,12	0
4	NA	A	3104	1/1	0.99	0.06	21,21,21,21	0
5	DMS	C	8403	4/4	0.99	0.10	10,12,16,19	0
5	DMS	C	8405	4/4	0.99	0.10	16,23,25,38	0
5	DMS	C	8412	4/4	0.99	0.06	18,31,67,100	0
4	NA	D	3101	1/1	0.99	0.04	16,16,16,16	0
3	MG	D	3001	1/1	0.99	0.02	15,15,15,15	0
5	DMS	D	8404	4/4	0.99	0.07	14,19,26,80	0
5	DMS	B	8404	4/4	0.99	0.07	16,20,30,40	0
5	DMS	A	8404	4/4	0.99	0.06	9,22,22,30	0
5	DMS	B	8403	4/4	0.99	0.12	12,12,18,21	0
5	DMS	A	8410	4/4	0.99	0.09	28,39,55,58	0
3	MG	A	3001	1/1	0.99	0.02	13,13,13,13	0
5	DMS	B	8401	4/4	0.99	0.07	13,13,21,22	0
5	DMS	C	8404	4/4	0.99	0.06	11,13,21,21	0
5	DMS	D	8411	4/4	0.99	0.05	11,27,27,100	0
5	DMS	A	8405	4/4	0.99	0.05	11,22,25,27	0
5	DMS	C	8401	4/4	0.99	0.08	10,11,21,30	0
5	DMS	C	8402	4/4	0.99	0.07	6,31,38,40	0
5	DMS	A	8403	4/4	0.99	0.09	15,20,21,23	0
4	NA	C	3102	1/1	0.99	0.04	16,16,16,16	0
5	DMS	D	8401	4/4	0.99	0.06	10,11,20,27	0
3	MG	C	3001	1/1	0.99	0.03	12,12,12,12	0
5	DMS	B	8411	4/4	0.99	0.05	15,17,32,100	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	DMS	A	8408	4/4	0.99	0.08	14,23,31,82	0
5	DMS	D	8405	4/4	0.99	0.08	18,23,42,100	0
5	DMS	D	8403	4/4	0.99	0.10	21,26,27,45	0
3	MG	B	3002	1/1	0.99	0.04	14,14,14,14	0
5	DMS	A	8401	4/4	0.99	0.09	6,13,16,19	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.