



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 03:06 PM BST

PDB ID : 3T0B  
Title : E. coli (LacZ) beta-galactosidase (S796T) IPTG complex  
Authors : Jancewicz, L.J.; Wheatley, R.W.; Sutendra, G.; Lee, M.; Fraser, M.; Huber, R.E.  
Deposited on : 2011-07-19  
Resolution : 2.40 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](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

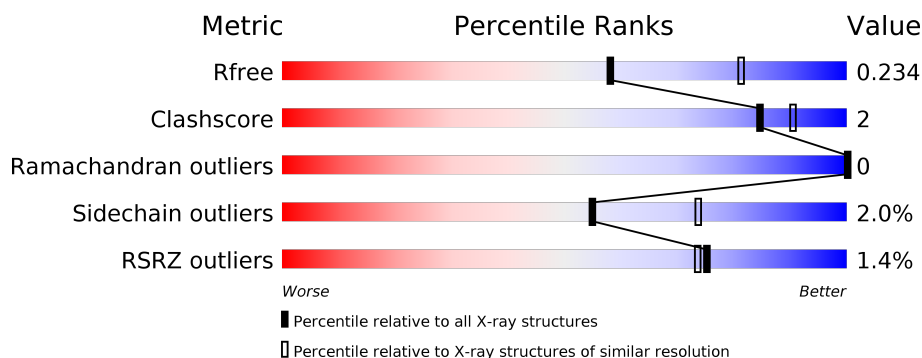
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	1052	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>7%</div> <div>.</div> </div> </div>
1	B	1052	<div> <div>%</div> <div> <div></div> <div>91%</div> <div>6%</div> <div>.</div> </div> </div>
1	C	1052	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>8%</div> <div>.</div> </div> </div>
1	D	1052	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>7%</div> <div>.</div> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 36706 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
			8151	5155	1444	1514	38			
1	B	1015	Total	C	N	O	S	0	0	0
			8158	5160	1445	1515	38			
1	C	1015	Total	C	N	O	S	0	0	0
			8158	5160	1445	1515	38			
1	D	1015	Total	C	N	O	S	0	0	0
			8158	5160	1445	1515	38			

There are 152 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-28	MET	-	expression tag	UNP P00722
A	-27	GLY	-	expression tag	UNP P00722
A	-26	GLY	-	expression tag	UNP P00722
A	-25	SER	-	expression tag	UNP P00722
A	-24	HIS	-	expression tag	UNP P00722
A	-23	HIS	-	expression tag	UNP P00722
A	-22	HIS	-	expression tag	UNP P00722
A	-21	HIS	-	expression tag	UNP P00722
A	-20	HIS	-	expression tag	UNP P00722
A	-19	HIS	-	expression tag	UNP P00722
A	-18	GLY	-	expression tag	UNP P00722
A	-17	MET	-	expression tag	UNP P00722
A	-16	ALA	-	expression tag	UNP P00722
A	-15	SER	-	expression tag	UNP P00722
A	-14	MET	-	expression tag	UNP P00722
A	-13	THR	-	expression tag	UNP P00722
A	-12	GLY	-	expression tag	UNP P00722
A	-11	GLY	-	expression tag	UNP P00722
A	-10	GLN	-	expression tag	UNP P00722
A	-9	GLN	-	expression tag	UNP P00722
A	-8	MET	-	expression tag	UNP P00722

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	expression tag	UNP P00722
A	-6	ARG	-	expression tag	UNP P00722
A	-5	ASP	-	expression tag	UNP P00722
A	-4	LEU	-	expression tag	UNP P00722
A	-3	TYR	-	expression tag	UNP P00722
A	-2	ASP	-	expression tag	UNP P00722
A	-1	ASP	-	expression tag	UNP P00722
A	0	ASP	-	expression tag	UNP P00722
A	1	ASP	-	expression tag	UNP P00722
A	2	LYS	-	expression tag	UNP P00722
A	3	ASP	-	expression tag	UNP P00722
A	4	PRO	-	expression tag	UNP P00722
A	5	MET	-	expression tag	UNP P00722
A	6	ILE	-	expression tag	UNP P00722
A	7	ASP	-	expression tag	UNP P00722
A	8	PRO	-	expression tag	UNP P00722
A	796	THR	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	THR	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	THR	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	THR	SER	engineered mutation	UNP P00722

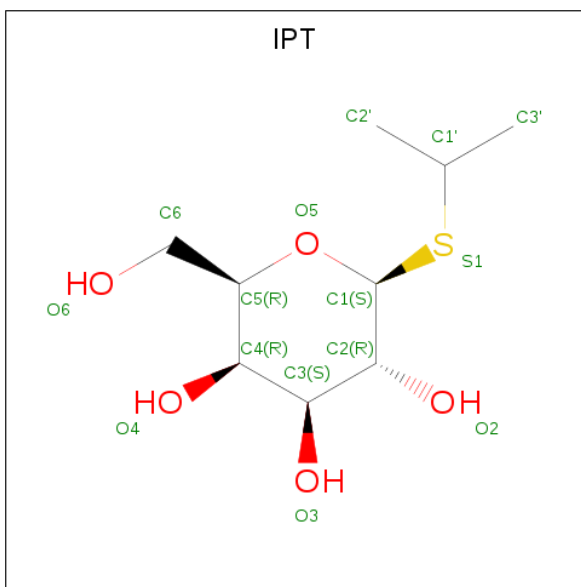
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

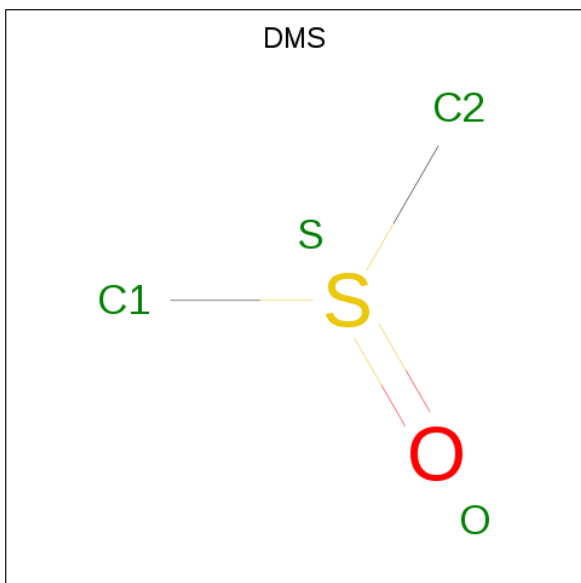
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	4	Total Na 4 4	0	0
3	A	3	Total Na 3 3	0	0
3	D	3	Total Na 3 3	0	0
3	C	4	Total Na 4 4	0	0

- Molecule 4 is 1-methylethyl 1-thio-beta-D-galactopyranoside (three-letter code: IPT) (formula: C<sub>9</sub>H<sub>18</sub>O<sub>5</sub>S).



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

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula:  $C_2H_6OS$ ).





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

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

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

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

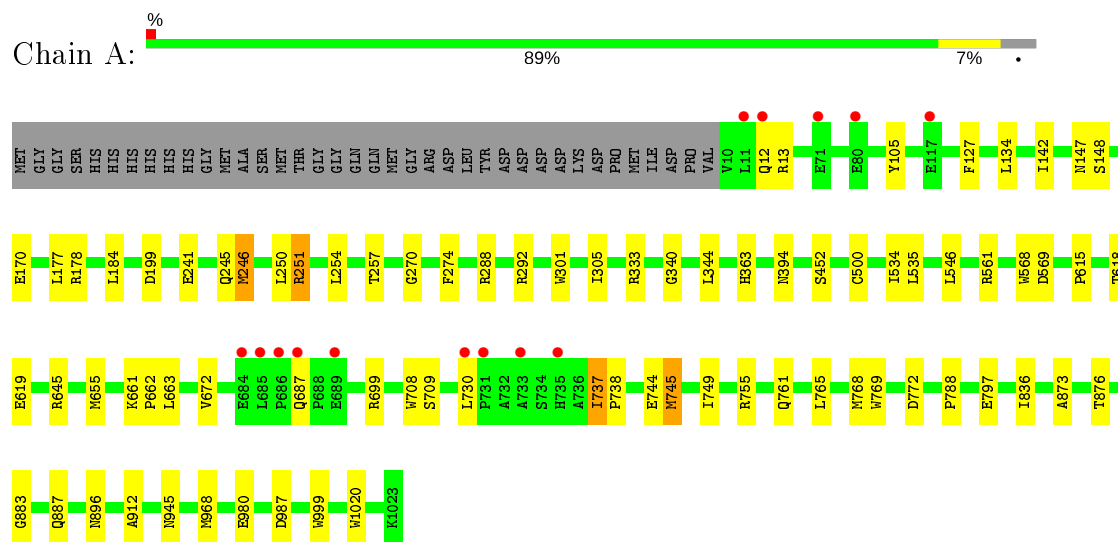
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	857	Total 857	O 857	0	0
6	B	992	Total 992	O 992	0	0
6	C	1017	Total 1017	O 1017	0	0
6	D	857	Total 857	O 857	0	0

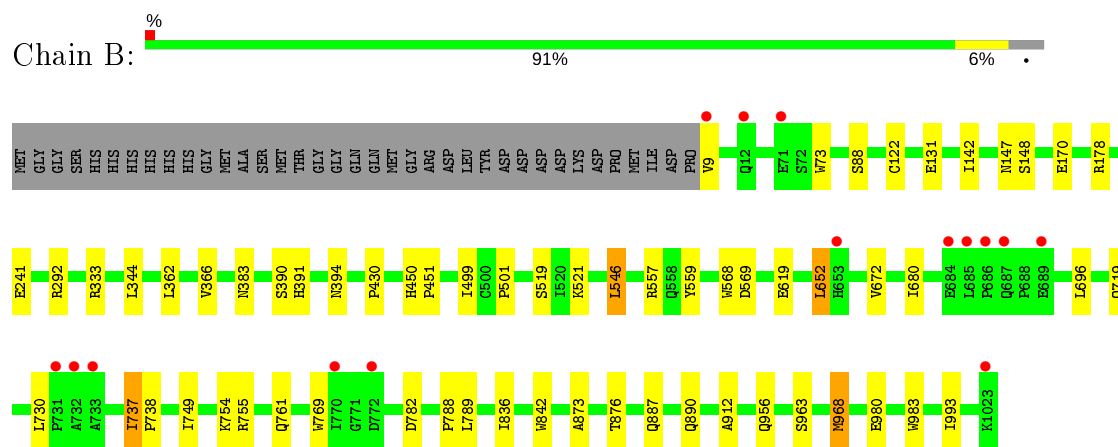
### 3 Residue-property plots [i](#)

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

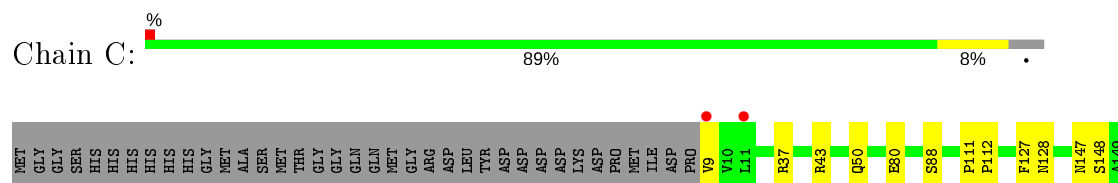
#### • Molecule 1: Beta-galactosidase

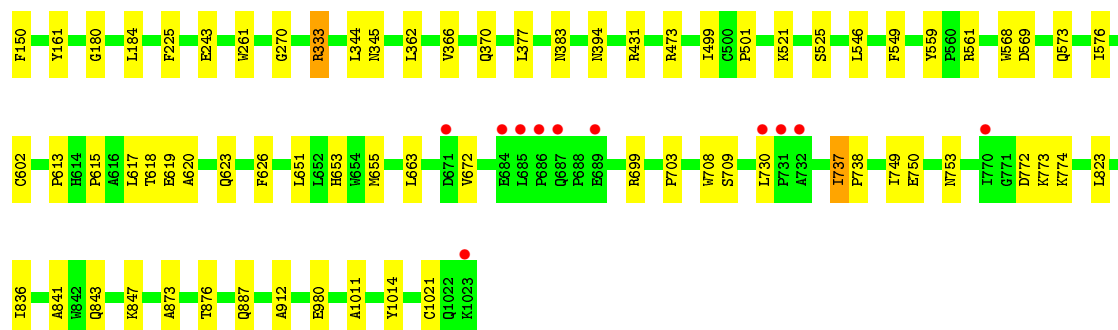


#### • Molecule 1: Beta-galactosidase

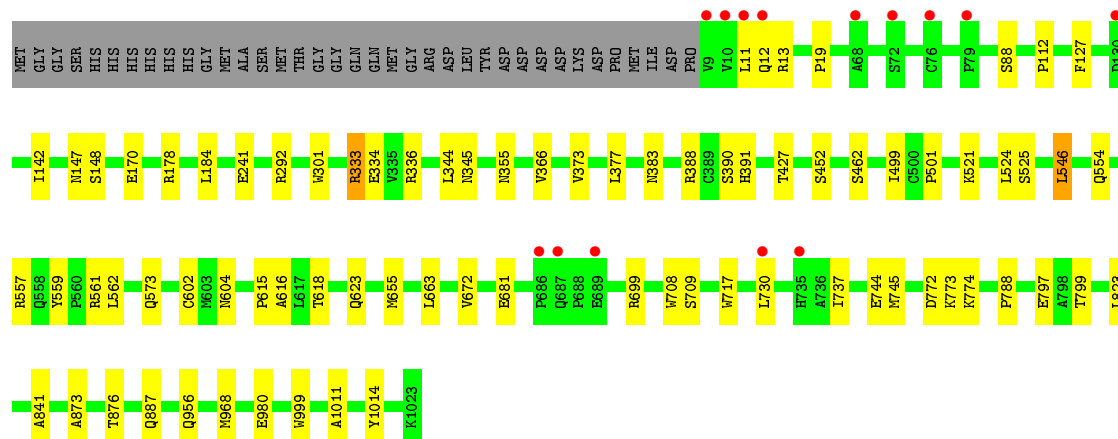
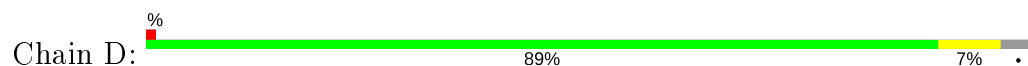


#### • Molecule 1: Beta-galactosidase





• Molecule 1: Beta-galactosidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	151.97Å 163.14Å 204.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	67.79 – 2.40 62.07 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (67.79-2.40) 98.0 (62.07-2.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.23 (at 2.40Å)	Xtrriage
Refinement program	REFMAC 5.5.0109, CNS	Depositor
R, $R_{free}$	0.165 , 0.233 0.167 , 0.234	Depositor DCC
$R_{free}$ test set	2826 reflections (1.46%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.4	Xtrriage
Anisotropy	0.081	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 45.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	36706	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	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 41.18 % 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 2.4517e-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 [i](#)

### 5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.40	0/8393	0.55	0/11450
1	B	0.41	0/8400	0.55	0/11460
1	C	0.42	0/8400	0.55	0/11460
1	D	0.39	0/8400	0.54	0/11460
All	All	0.40	0/33593	0.55	0/45830

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	8151	0	7746	40	0
1	B	8158	0	7755	29	0
1	C	8158	0	7755	45	0
1	D	8158	0	7755	36	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	3	0	0	0	0

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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	15	0	17	1	0
4	B	15	0	17	0	0
4	C	15	0	17	0	0
4	D	15	0	17	1	0
5	A	64	0	96	2	0
5	B	72	0	108	3	0
5	C	76	0	114	8	0
5	D	64	0	96	4	0
6	A	857	0	0	2	0
6	B	992	0	0	2	0
6	C	1017	0	0	5	0
6	D	857	0	0	2	0
All	All	36706	0	31493	148	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 (148) 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:383:ASN:HA	5:C:8002:DMS:H21	1.59	0.82
1:A:245:GLN:HG2	1:A:288:ARG:HG2	1.66	0.76
1:A:270:GLY:HA3	5:A:8005:DMS:H21	1.73	0.70
1:C:887:GLN:NE2	1:C:980:GLU:O	2.24	0.70
1:A:241:GLU:HG3	1:A:292:ARG:HG2	1.81	0.62
1:B:383:ASN:HA	5:B:8003:DMS:H11	1.81	0.61
1:B:142:ILE:HG12	1:B:170:GLU:HG2	1.82	0.61
1:A:699:ARG:NH1	6:A:4812:HOH:O	2.33	0.61
1:C:499:ILE:HG22	1:C:501:PRO:HD3	1.83	0.61
1:D:127:PHE:CE2	1:D:184:LEU:HG	2.37	0.59
1:B:131:GLU:HG2	6:B:4785:HOH:O	2.02	0.58
1:A:246:MET:CE	1:A:254:LEU:HD13	2.32	0.58
1:D:499:ILE:HG22	1:D:501:PRO:HD3	1.85	0.57
1:D:142:ILE:HG12	1:D:170:GLU:HG2	1.87	0.57
1:A:246:MET:HG2	1:A:274:PHE:CE2	2.40	0.56
1:D:887:GLN:NE2	1:D:980:GLU:O	2.32	0.56
1:B:887:GLN:NE2	1:B:980:GLU:O	2.36	0.55
1:A:788:PRO:HD2	1:A:968:MET:HB2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:377:LEU:HD22	1:C:708:TRP:HA	1.89	0.55
1:A:270:GLY:HA3	5:A:8005:DMS:C2	2.37	0.55
1:A:142:ILE:HG12	1:A:170:GLU:HG2	1.91	0.53
1:C:88:SER:HA	1:C:366:VAL:HG21	1.90	0.53
1:B:499:ILE:HG22	1:B:501:PRO:HD3	1.91	0.53
1:B:737:ILE:HG12	1:B:738:PRO:HD2	1.91	0.53
1:C:623:GLN:HA	5:C:8001:DMS:O	2.08	0.53
1:D:355:ASN:OD1	1:D:388:ARG:HD3	2.09	0.53
1:A:887:GLN:NE2	1:A:980:GLU:O	2.39	0.52
1:A:765:LEU:HD21	1:A:768:MET:HE2	1.90	0.52
1:C:127:PHE:CE2	1:C:184:LEU:HG	2.45	0.52
1:D:788:PRO:HD2	1:D:968:MET:HB2	1.92	0.52
1:B:652:LEU:HD22	1:B:680:ILE:HD13	1.93	0.51
1:A:615:PRO:O	1:A:618:THR:HG22	2.10	0.51
1:D:127:PHE:HE2	1:D:184:LEU:HG	1.74	0.51
1:D:427:THR:HG21	1:D:462:SER:HB3	1.92	0.51
1:A:619:GLU:HA	1:A:912:ALA:HB2	1.92	0.50
1:C:431:ARG:HD3	6:C:4676:HOH:O	2.10	0.50
1:C:576:ILE:HD12	5:C:8008:DMS:H22	1.93	0.50
1:C:699:ARG:NH1	5:C:8011:DMS:H13	2.27	0.50
1:D:377:LEU:HD22	1:D:708:TRP:HA	1.93	0.50
1:A:708:TRP:CE3	1:A:709:SER:HB3	2.46	0.50
1:D:334:GLU:OE1	1:D:336:ARG:NH1	2.44	0.49
1:A:127:PHE:HE2	1:A:184:LEU:HG	1.78	0.49
1:B:557:ARG:NH1	5:B:8002:DMS:O	2.46	0.49
1:C:753:ASN:HB3	6:C:4002:HOH:O	2.13	0.49
1:A:12:GLN:O	1:A:13:ARG:HB2	2.13	0.48
1:C:333:ARG:HA	1:C:345:ASN:OD1	2.13	0.48
1:B:521:LYS:HG2	1:B:559:TYR:CZ	2.49	0.48
1:C:383:ASN:HA	5:C:8002:DMS:C2	2.38	0.48
1:D:373:VAL:O	1:D:377:LEU:HG	2.13	0.48
1:D:383:ASN:HA	5:D:8003:DMS:H11	1.96	0.48
1:D:521:LYS:HG2	1:D:559:TYR:CZ	2.49	0.48
1:C:651:LEU:HD23	1:C:703:PRO:HG3	1.95	0.48
1:A:873:ALA:O	1:A:876:THR:HG22	2.14	0.47
1:B:789:LEU:HD11	1:B:993:ILE:HG22	1.96	0.47
1:C:549:PHE:CE2	1:C:620:ALA:HA	2.48	0.47
1:C:651:LEU:HD11	1:C:653:HIS:CE1	2.49	0.47
1:C:1011:ALA:HB3	1:C:1014:TYR:CZ	2.49	0.47
1:D:999:TRP:CD2	4:D:2001:IPT:H3'3	2.49	0.47
1:B:430:PRO:HD3	5:B:8010:DMS:H22	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:SER:HA	1:B:366:VAL:HG21	1.96	0.47
1:C:576:ILE:HD12	5:C:8008:DMS:C2	2.45	0.47
1:A:147:ASN:HA	1:A:148:SER:HA	1.70	0.47
1:A:105:TYR:CE2	1:A:199:ASP:HB2	2.50	0.47
1:B:788:PRO:HD2	1:B:968:MET:HB2	1.97	0.46
1:C:147:ASN:HA	1:C:148:SER:HA	1.66	0.46
1:D:573:GLN:HB2	1:D:602:CYS:O	2.16	0.46
1:C:225:PHE:HA	1:C:243:GLU:O	2.15	0.46
1:C:749:ILE:HD11	1:C:836:ILE:HD11	1.96	0.46
1:C:37:ARG:HG2	1:C:50:GLN:NE2	2.30	0.46
1:B:9:VAL:HG22	1:C:9:VAL:HG22	1.98	0.46
1:D:147:ASN:HA	1:D:148:SER:HA	1.64	0.46
1:A:127:PHE:CE2	1:A:184:LEU:HG	2.51	0.46
1:C:626:PHE:HE1	5:C:8002:DMS:H21	1.80	0.46
1:A:999:TRP:CD2	4:A:2001:IPT:H3'3	2.50	0.45
1:D:1011:ALA:HB3	1:D:1014:TYR:CZ	2.52	0.45
1:A:246:MET:HE1	1:A:254:LEU:HD13	1.96	0.45
1:B:390:SER:HA	1:B:391:HIS:HA	1.79	0.45
1:C:873:ALA:O	1:C:876:THR:HG22	2.16	0.45
1:D:557:ARG:NH1	5:D:8002:DMS:O	2.49	0.45
1:D:873:ALA:O	1:D:876:THR:HG22	2.16	0.45
1:D:19:PRO:HD3	1:D:112:PRO:HB3	1.99	0.44
1:D:623:GLN:HA	5:D:8002:DMS:O	2.17	0.44
1:D:390:SER:HA	1:D:391:HIS:HA	1.76	0.44
1:C:473:ARG:NH1	6:C:4634:HOH:O	2.42	0.44
1:A:737:ILE:HA	1:A:738:PRO:HD3	1.88	0.44
1:B:546:LEU:HA	6:B:4134:HOH:O	2.18	0.44
1:C:270:GLY:HA3	5:C:8004:DMS:O	2.17	0.44
1:D:615:PRO:O	1:D:618:THR:HG22	2.17	0.44
1:B:696:LEU:O	1:B:719:GLN:HA	2.17	0.44
1:C:561:ARG:HD3	1:D:525:SER:O	2.17	0.44
1:C:615:PRO:O	1:C:618:THR:HG22	2.17	0.44
1:C:750:GLU:HB2	6:C:4649:HOH:O	2.18	0.44
1:B:873:ALA:O	1:B:876:THR:HG22	2.19	0.43
1:A:250:LEU:C	1:A:251:ARG:HG2	2.39	0.43
1:C:619:GLU:HA	1:C:912:ALA:HB2	2.01	0.43
1:D:88:SER:HA	1:D:366:VAL:HG21	2.01	0.43
1:A:883:GLY:HA3	1:A:987:ASP:HA	1.99	0.43
1:D:524:LEU:HD11	1:D:562:LEU:HG	2.01	0.43
1:C:43:ARG:HD2	1:C:261:TRP:CE3	2.54	0.43
1:B:754:LYS:HA	1:B:769:TRP:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:568:TRP:CD2	1:A:569:ASP:HB3	2.54	0.43
1:B:568:TRP:CD2	1:B:569:ASP:HB3	2.54	0.43
1:B:755:ARG:HB3	1:B:769:TRP:HB2	1.99	0.43
1:C:521:LYS:HG2	1:C:559:TYR:CZ	2.54	0.43
1:D:12:GLN:O	1:D:13:ARG:HB2	2.18	0.43
1:A:301:TRP:CH2	1:A:452:SER:HA	2.55	0.42
1:B:147:ASN:HA	1:B:148:SER:HA	1.64	0.42
1:C:573:GLN:HB2	1:C:602:CYS:O	2.19	0.42
1:D:823:LEU:HD11	1:D:841:ALA:HB2	2.01	0.42
1:A:305:ILE:HD11	1:A:645:ARG:HB3	2.02	0.42
1:C:1021:CYS:HB2	6:C:4972:HOH:O	2.19	0.42
1:C:568:TRP:CD2	1:C:569:ASP:HB3	2.54	0.42
1:C:128:ASN:HA	1:C:180:GLY:O	2.19	0.42
5:D:8006:DMS:H12	6:D:4368:HOH:O	2.18	0.42
1:C:737:ILE:HD13	1:C:738:PRO:O	2.19	0.42
1:D:699:ARG:HG2	1:D:717:TRP:HB3	2.02	0.42
1:B:782:ASP:HB2	1:B:842:TRP:CZ2	2.55	0.42
1:C:111:PRO:HA	1:C:112:PRO:HA	1.85	0.42
1:A:896:ASN:HB3	1:A:945:ASN:HB2	2.02	0.42
1:D:301:TRP:CH2	1:D:452:SER:HA	2.54	0.42
1:A:749:ILE:HD11	1:A:836:ILE:HD11	2.01	0.41
1:C:708:TRP:CE3	1:C:709:SER:HB3	2.55	0.41
1:B:963:SER:HB3	1:B:983:TRP:CE2	2.54	0.41
1:D:546:LEU:HD22	1:D:616:ALA:HB1	2.01	0.41
1:D:333:ARG:HA	1:D:345:ASN:OD1	2.21	0.41
1:D:708:TRP:CE3	1:D:709:SER:HB3	2.55	0.41
1:C:150:PHE:O	1:C:161:TYR:HA	2.21	0.41
1:A:745:MET:HG3	1:A:761:GLN:OE1	2.21	0.41
1:B:241:GLU:HG3	1:B:292:ARG:HG2	2.02	0.41
1:B:450:HIS:HA	1:B:451:PRO:HD2	1.94	0.41
1:C:525:SER:O	1:D:561:ARG:HD3	2.20	0.41
1:A:768:MET:HE3	1:A:1020:TRP:CZ2	2.56	0.41
1:A:363:HIS:HD2	6:A:4479:HOH:O	2.03	0.41
1:A:134:LEU:HD21	1:A:177:LEU:HB2	2.03	0.41
1:B:619:GLU:HA	1:B:912:ALA:HB2	2.02	0.41
1:C:613:PRO:HB3	1:C:617:LEU:HD23	2.03	0.40
1:D:241:GLU:HG3	1:D:292:ARG:HG2	2.03	0.40
1:B:73:TRP:CE2	1:B:122:CYS:HB3	2.56	0.40
1:B:749:ILE:HD11	1:B:836:ILE:HD11	2.02	0.40
1:C:843:GLN:HA	1:C:847:LYS:O	2.22	0.40
1:A:340:GLY:O	1:A:561:ARG:HG2	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:CYS:HA	1:A:534:ILE:O	2.21	0.40
1:A:755:ARG:HB3	1:A:769:TRP:HE3	1.87	0.40
1:A:246:MET:SD	1:A:250:LEU:HD23	2.61	0.40
1:A:257:THR:HA	1:A:270:GLY:O	2.20	0.40
1:A:661:LYS:HA	1:A:662:PRO:HD3	1.96	0.40
1:C:823:LEU:HD11	1:C:841:ALA:HB2	2.02	0.40
1:D:799:THR:HG22	6:D:4759:HOH:O	2.22	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%)	976 (96%)	36 (4%)	0	100	100
1	B	1013/1052 (96%)	976 (96%)	37 (4%)	0	100	100
1	C	1013/1052 (96%)	979 (97%)	34 (3%)	0	100	100
1	D	1013/1052 (96%)	982 (97%)	31 (3%)	0	100	100
All	All	4051/4208 (96%)	3913 (97%)	138 (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	867/898 (96%)	849 (98%)	18 (2%)	53	72
1	B	868/898 (97%)	853 (98%)	15 (2%)	60	78
1	C	868/898 (97%)	853 (98%)	15 (2%)	60	78
1	D	868/898 (97%)	848 (98%)	20 (2%)	50	70
All	All	3471/3592 (97%)	3403 (98%)	68 (2%)	55	74

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	178	ARG
1	A	246	MET
1	A	251	ARG
1	A	333	ARG
1	A	344	LEU
1	A	394	ASN
1	A	535	LEU
1	A	546	LEU
1	A	655	MET
1	A	663	LEU
1	A	672	VAL
1	A	687	GLN
1	A	730	LEU
1	A	737	ILE
1	A	744	GLU
1	A	745	MET
1	A	772	ASP
1	A	797	GLU
1	B	178	ARG
1	B	333	ARG
1	B	344	LEU
1	B	362	LEU
1	B	394	ASN
1	B	519	SER
1	B	546	LEU
1	B	652	LEU
1	B	672	VAL
1	B	730	LEU
1	B	737	ILE
1	B	761	GLN
1	B	890	GLN
1	B	956	GLN
1	B	968	MET

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Mol	Chain	Res	Type
1	C	80	GLU
1	C	333	ARG
1	C	344	LEU
1	C	362	LEU
1	C	370	GLN
1	C	394	ASN
1	C	546	LEU
1	C	655	MET
1	C	663	LEU
1	C	672	VAL
1	C	730	LEU
1	C	737	ILE
1	C	772	ASP
1	C	773	LYS
1	C	774	LYS
1	D	11	LEU
1	D	178	ARG
1	D	333	ARG
1	D	344	LEU
1	D	546	LEU
1	D	554	GLN
1	D	604	ASN
1	D	655	MET
1	D	663	LEU
1	D	672	VAL
1	D	681	GLU
1	D	730	LEU
1	D	737	ILE
1	D	744	GLU
1	D	745	MET
1	D	772	ASP
1	D	773	LYS
1	D	774	LYS
1	D	797	GLU
1	D	956	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	363	HIS
1	A	370	GLN
1	A	775	GLN

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Mol	Chain	Res	Type
1	C	50	GLN
1	C	510	GLN
1	C	739	HIS
1	C	1017	GLN
1	D	757	GLN
1	D	863	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 95 ligands modelled in this entry, 22 are monoatomic - leaving 73 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	D	8002	-	3,3,3	2.54	1 (33%)	3,3,3	0.29	0
4	IPT	C	2001	3	14,15,15	0.52	0	18,21,21	1.35	2 (11%)
5	DMS	B	8002	-	3,3,3	2.56	1 (33%)	3,3,3	0.52	0
5	DMS	D	8016	-	3,3,3	2.64	1 (33%)	3,3,3	0.50	0
5	DMS	C	8001	-	3,3,3	2.56	1 (33%)	3,3,3	0.25	0
5	DMS	C	8008	-	3,3,3	2.62	1 (33%)	3,3,3	0.64	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	DMS	D	8001	-	3,3,3	2.63	1 (33%)	3,3,3	0.86	0
5	DMS	B	8005	-	3,3,3	2.63	1 (33%)	3,3,3	0.47	0
5	DMS	A	8001	-	3,3,3	2.63	1 (33%)	3,3,3	0.60	0
5	DMS	C	8002	-	3,3,3	2.47	1 (33%)	3,3,3	0.56	0
5	DMS	A	8011	-	3,3,3	2.63	1 (33%)	3,3,3	0.56	0
5	DMS	A	8005	-	3,3,3	2.61	1 (33%)	3,3,3	0.28	0
5	DMS	B	8010	-	3,3,3	2.64	1 (33%)	3,3,3	0.67	0
5	DMS	A	8015	-	3,3,3	2.65	1 (33%)	3,3,3	0.61	0
5	DMS	D	8004	-	3,3,3	2.60	1 (33%)	3,3,3	0.54	0
5	DMS	D	8014	-	3,3,3	2.67	1 (33%)	3,3,3	0.56	0
4	IPT	D	2001	3	14,15,15	0.50	0	18,21,21	1.43	1 (5%)
5	DMS	A	8014	-	3,3,3	2.68	1 (33%)	3,3,3	0.52	0
5	DMS	D	8003	-	3,3,3	2.63	1 (33%)	3,3,3	0.54	0
5	DMS	C	8003	-	3,3,3	2.67	1 (33%)	3,3,3	0.57	0
5	DMS	C	8018	-	3,3,3	2.63	1 (33%)	3,3,3	0.49	0
5	DMS	B	8009	-	3,3,3	2.59	1 (33%)	3,3,3	0.61	0
5	DMS	C	8014	-	3,3,3	2.64	1 (33%)	3,3,3	0.49	0
5	DMS	C	8017	-	3,3,3	2.61	1 (33%)	3,3,3	0.61	0
5	DMS	B	8015	-	3,3,3	2.68	1 (33%)	3,3,3	0.49	0
5	DMS	C	8005	-	3,3,3	2.62	1 (33%)	3,3,3	0.47	0
5	DMS	C	8009	-	3,3,3	2.65	1 (33%)	3,3,3	0.67	0
5	DMS	B	8001	-	3,3,3	2.53	1 (33%)	3,3,3	0.76	0
5	DMS	C	8006	-	3,3,3	2.70	1 (33%)	3,3,3	0.68	0
5	DMS	B	8017	-	3,3,3	2.68	1 (33%)	3,3,3	0.57	0
5	DMS	A	8009	-	3,3,3	2.66	1 (33%)	3,3,3	0.53	0
4	IPT	A	2001	3	14,15,15	0.51	0	18,21,21	1.32	1 (5%)
5	DMS	A	8003	-	3,3,3	2.66	1 (33%)	3,3,3	0.61	0
5	DMS	D	8013	-	3,3,3	2.67	1 (33%)	3,3,3	0.51	0
5	DMS	B	8014	-	3,3,3	2.62	1 (33%)	3,3,3	0.58	0
5	DMS	D	8015	-	3,3,3	2.65	1 (33%)	3,3,3	0.53	0
5	DMS	D	8009	-	3,3,3	2.64	1 (33%)	3,3,3	0.56	0
5	DMS	D	8011	-	3,3,3	2.66	1 (33%)	3,3,3	0.73	0
5	DMS	B	8011	-	3,3,3	2.62	1 (33%)	3,3,3	0.53	0
5	DMS	D	8005	-	3,3,3	2.59	1 (33%)	3,3,3	0.53	0
5	DMS	C	8012	-	3,3,3	2.62	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	B	8012	-	3,3,3	2.64	1 (33%)	3,3,3	0.47	0
5	DMS	D	8012	-	3,3,3	2.63	1 (33%)	3,3,3	0.48	0
4	IPT	B	2001	3	14,15,15	0.52	0	18,21,21	1.74	2 (11%)
5	DMS	B	8016	-	3,3,3	2.63	1 (33%)	3,3,3	0.45	0
5	DMS	A	8012	-	3,3,3	2.64	1 (33%)	3,3,3	0.46	0
5	DMS	B	8018	-	3,3,3	2.64	1 (33%)	3,3,3	0.66	0
5	DMS	A	8002	-	3,3,3	2.61	1 (33%)	3,3,3	0.53	0
5	DMS	A	8016	-	3,3,3	2.66	1 (33%)	3,3,3	0.55	0
5	DMS	A	8006	-	3,3,3	2.70	1 (33%)	3,3,3	0.54	0
5	DMS	C	8011	-	3,3,3	2.65	1 (33%)	3,3,3	0.63	0
5	DMS	D	8007	-	3,3,3	2.58	1 (33%)	3,3,3	0.38	0
5	DMS	A	8007	-	3,3,3	2.68	1 (33%)	3,3,3	0.57	0
5	DMS	D	8010	-	3,3,3	2.66	1 (33%)	3,3,3	0.51	0
5	DMS	C	8010	-	3,3,3	2.61	1 (33%)	3,3,3	0.62	0
5	DMS	C	8013	-	3,3,3	2.66	1 (33%)	3,3,3	0.57	0
5	DMS	C	8007	-	3,3,3	2.65	1 (33%)	3,3,3	0.49	0
5	DMS	A	8004	-	3,3,3	2.66	1 (33%)	3,3,3	0.56	0
5	DMS	C	8019	-	3,3,3	2.56	1 (33%)	3,3,3	0.44	0
5	DMS	C	8004	-	3,3,3	2.63	1 (33%)	3,3,3	0.53	0
5	DMS	A	8008	-	3,3,3	2.58	1 (33%)	3,3,3	0.35	0
5	DMS	B	8006	-	3,3,3	2.67	1 (33%)	3,3,3	0.72	0
5	DMS	C	8016	-	3,3,3	2.60	1 (33%)	3,3,3	0.47	0
5	DMS	B	8013	-	3,3,3	2.63	1 (33%)	3,3,3	0.55	0
5	DMS	D	8008	-	3,3,3	2.58	1 (33%)	3,3,3	0.52	0
5	DMS	C	8015	-	3,3,3	2.69	1 (33%)	3,3,3	0.50	0
5	DMS	B	8004	-	3,3,3	2.57	1 (33%)	3,3,3	0.59	0
5	DMS	B	8008	-	3,3,3	2.54	1 (33%)	3,3,3	0.60	0
5	DMS	B	8003	-	3,3,3	2.67	1 (33%)	3,3,3	0.63	0
5	DMS	A	8010	-	3,3,3	2.62	1 (33%)	3,3,3	0.57	0
5	DMS	B	8007	-	3,3,3	2.65	1 (33%)	3,3,3	0.58	0
5	DMS	A	8013	-	3,3,3	2.63	1 (33%)	3,3,3	0.32	0
5	DMS	D	8006	-	3,3,3	2.67	1 (33%)	3,3,3	0.58	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.  
'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	IPT	C	2001	3	-	4/6/26/26	0/1/1/1
4	IPT	D	2001	3	-	3/6/26/26	0/1/1/1
4	IPT	A	2001	3	-	3/6/26/26	0/1/1/1
4	IPT	B	2001	3	-	4/6/26/26	0/1/1/1

All (69) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	8006	DMS	O-S	4.54	1.80	1.50
5	C	8006	DMS	O-S	4.52	1.80	1.50
5	C	8003	DMS	O-S	4.52	1.80	1.50
5	C	8015	DMS	O-S	4.52	1.80	1.50
5	B	8003	DMS	O-S	4.50	1.80	1.50
5	B	8017	DMS	O-S	4.49	1.80	1.50
5	B	8015	DMS	O-S	4.49	1.80	1.50
5	A	8014	DMS	O-S	4.49	1.80	1.50
5	D	8006	DMS	O-S	4.49	1.80	1.50
5	D	8014	DMS	O-S	4.48	1.80	1.50
5	B	8006	DMS	O-S	4.48	1.80	1.50
5	A	8007	DMS	O-S	4.48	1.80	1.50
5	A	8003	DMS	O-S	4.48	1.80	1.50
5	D	8013	DMS	O-S	4.47	1.80	1.50
5	A	8004	DMS	O-S	4.47	1.80	1.50
5	A	8016	DMS	O-S	4.47	1.80	1.50
5	D	8011	DMS	O-S	4.46	1.80	1.50
5	D	8010	DMS	O-S	4.46	1.80	1.50
5	C	8009	DMS	O-S	4.45	1.80	1.50
5	C	8013	DMS	O-S	4.45	1.80	1.50
5	A	8009	DMS	O-S	4.45	1.80	1.50
5	B	8013	DMS	O-S	4.45	1.80	1.50
5	D	8015	DMS	O-S	4.44	1.80	1.50
5	A	8012	DMS	O-S	4.44	1.80	1.50
5	C	8011	DMS	O-S	4.44	1.80	1.50
5	A	8015	DMS	O-S	4.44	1.80	1.50
5	D	8016	DMS	O-S	4.44	1.80	1.50
5	B	8007	DMS	O-S	4.44	1.80	1.50
5	B	8012	DMS	O-S	4.43	1.80	1.50
5	C	8007	DMS	O-S	4.43	1.80	1.50
5	B	8010	DMS	O-S	4.43	1.80	1.50
5	B	8018	DMS	O-S	4.43	1.80	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	8014	DMS	O-S	4.43	1.80	1.50
5	D	8012	DMS	O-S	4.43	1.80	1.50
5	D	8003	DMS	O-S	4.42	1.80	1.50
5	C	8008	DMS	O-S	4.42	1.80	1.50
5	D	8009	DMS	O-S	4.42	1.80	1.50
5	A	8013	DMS	O-S	4.42	1.80	1.50
5	A	8001	DMS	O-S	4.42	1.80	1.50
5	A	8005	DMS	O-S	4.42	1.80	1.50
5	C	8005	DMS	O-S	4.41	1.80	1.50
5	B	8016	DMS	O-S	4.40	1.80	1.50
5	C	8018	DMS	O-S	4.40	1.80	1.50
5	A	8011	DMS	O-S	4.40	1.80	1.50
5	B	8005	DMS	O-S	4.40	1.79	1.50
5	D	8001	DMS	O-S	4.40	1.79	1.50
5	C	8004	DMS	O-S	4.40	1.79	1.50
5	B	8011	DMS	O-S	4.39	1.79	1.50
5	A	8002	DMS	O-S	4.39	1.79	1.50
5	C	8012	DMS	O-S	4.38	1.79	1.50
5	C	8010	DMS	O-S	4.38	1.79	1.50
5	A	8010	DMS	O-S	4.38	1.79	1.50
5	C	8016	DMS	O-S	4.37	1.79	1.50
5	C	8017	DMS	O-S	4.37	1.79	1.50
5	B	8014	DMS	O-S	4.37	1.79	1.50
5	D	8004	DMS	O-S	4.36	1.79	1.50
5	B	8009	DMS	O-S	4.35	1.79	1.50
5	D	8005	DMS	O-S	4.34	1.79	1.50
5	A	8008	DMS	O-S	4.33	1.79	1.50
5	D	8008	DMS	O-S	4.32	1.79	1.50
5	D	8007	DMS	O-S	4.32	1.79	1.50
5	B	8004	DMS	O-S	4.32	1.79	1.50
5	B	8002	DMS	O-S	4.31	1.79	1.50
5	C	8001	DMS	O-S	4.31	1.79	1.50
5	C	8019	DMS	O-S	4.30	1.79	1.50
5	B	8008	DMS	O-S	4.28	1.79	1.50
5	D	8002	DMS	O-S	4.26	1.79	1.50
5	B	8001	DMS	O-S	4.23	1.78	1.50
5	C	8002	DMS	O-S	4.19	1.78	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	2001	IPT	C1-S1-C1'	4.86	115.14	100.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	2001	IPT	C1-S1-C1'	4.74	114.76	100.26
4	A	2001	IPT	C1-S1-C1'	4.72	114.70	100.26
4	B	2001	IPT	C2-C1-S1	-4.13	104.93	111.30
4	C	2001	IPT	C1-S1-C1'	3.99	112.48	100.26
4	C	2001	IPT	C2-C1-S1	-2.50	107.45	111.30

There are no chirality outliers.

All (14) torsion outliers are listed below:

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

There are no ring outliers.

14 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	8002	DMS	2	0
5	B	8002	DMS	1	0
5	C	8001	DMS	1	0
5	C	8008	DMS	2	0
5	C	8002	DMS	3	0
5	A	8005	DMS	2	0
5	B	8010	DMS	1	0
4	D	2001	IPT	1	0
5	D	8003	DMS	1	0
4	A	2001	IPT	1	0
5	C	8011	DMS	1	0
5	C	8004	DMS	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	8003	DMS	1	0
5	D	8006	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, 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	1014/1052 (96%)	-0.47	14 (1%) 75 73	7, 21, 44, 79	0
1	B	1015/1052 (96%)	-0.55	15 (1%) 73 72	7, 16, 40, 78	0
1	C	1015/1052 (96%)	-0.57	13 (1%) 77 75	6, 16, 38, 82	0
1	D	1015/1052 (96%)	-0.42	14 (1%) 75 73	10, 21, 45, 85	0
All	All	4059/4208 (96%)	-0.50	56 (1%) 75 73	6, 18, 42, 85	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	689	GLU	5.9
1	B	731	PRO	5.7
1	B	689	GLU	4.8
1	A	686	PRO	4.6
1	A	689	GLU	4.6
1	B	12	GLN	4.4
1	A	12	GLN	3.9
1	C	1023	LYS	3.8
1	D	687	GLN	3.8
1	B	770	ILE	3.6
1	D	686	PRO	3.5
1	D	11	LEU	3.5
1	B	684	GLU	3.4
1	B	732	ALA	3.4
1	B	686	PRO	3.4
1	D	9	VAL	3.4
1	C	686	PRO	3.3
1	D	130	ASP	3.2
1	C	689	GLU	3.1
1	C	687	GLN	3.1
1	A	730	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	730	LEU	3.0
1	B	1023	LYS	2.9
1	A	735	HIS	2.9
1	B	733	ALA	2.8
1	A	684	GLU	2.8
1	C	684	GLU	2.8
1	B	653	HIS	2.7
1	D	735	HIS	2.7
1	A	685	LEU	2.7
1	C	770	ILE	2.6
1	A	11	LEU	2.6
1	A	71	GLU	2.6
1	A	687	GLN	2.6
1	B	71	GLU	2.5
1	D	12	GLN	2.5
1	A	80	GLU	2.5
1	D	76	CYS	2.5
1	B	9	VAL	2.4
1	A	731	PRO	2.4
1	D	730	LEU	2.4
1	D	10	VAL	2.4
1	C	732	ALA	2.4
1	B	772	ASP	2.3
1	D	68	ALA	2.3
1	C	9	VAL	2.3
1	D	79	PRO	2.2
1	B	687	GLN	2.2
1	C	11	LEU	2.2
1	C	731	PRO	2.2
1	D	72	SER	2.1
1	B	685	LEU	2.1
1	C	685	LEU	2.1
1	A	117	GLU	2.1
1	C	671	ASP	2.1
1	A	733	ALA	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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



## 6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

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	A	8014	4/4	0.77	0.21	74,75,75,78	0
5	DMS	C	8014	4/4	0.80	0.41	61,64,65,67	0
5	DMS	C	8015	4/4	0.81	0.18	66,67,68,71	0
5	DMS	C	8006	4/4	0.84	0.23	55,56,59,61	0
5	DMS	A	8006	4/4	0.85	0.23	54,57,59,62	0
2	MG	A	3002	1/1	0.85	0.10	22,22,22,22	0
5	DMS	B	8012	4/4	0.85	0.40	41,45,48,49	0
5	DMS	A	8011	4/4	0.86	0.26	67,69,70,70	0
5	DMS	C	8011	4/4	0.87	0.24	58,58,61,62	0
5	DMS	D	8002	4/4	0.87	0.28	39,40,40,47	0
5	DMS	B	8007	4/4	0.87	0.27	52,56,57,57	0
2	MG	D	3002	1/1	0.87	0.19	24,24,24,24	0
5	DMS	D	8006	4/4	0.88	0.28	64,66,67,68	0
5	DMS	C	8010	4/4	0.88	0.25	48,50,51,55	0
5	DMS	A	8012	4/4	0.88	0.25	63,64,65,67	0
5	DMS	C	8001	4/4	0.89	0.21	27,32,35,39	0
5	DMS	D	8012	4/4	0.89	0.18	55,55,56,60	0
5	DMS	C	8007	4/4	0.89	0.27	39,44,47,47	0
3	NA	B	3104	1/1	0.90	0.19	36,36,36,36	0
5	DMS	A	8013	4/4	0.90	0.20	54,57,58,60	0
5	DMS	D	8010	4/4	0.90	0.29	65,67,68,70	0
3	NA	A	3103	1/1	0.91	0.16	31,31,31,31	0
5	DMS	D	8013	4/4	0.91	0.17	60,60,62,63	0
5	DMS	D	8014	4/4	0.91	0.17	49,50,51,53	0
5	DMS	C	8018	4/4	0.91	0.21	73,73,73,74	0
5	DMS	B	8006	4/4	0.91	0.23	47,49,50,51	0
5	DMS	D	8007	4/4	0.91	0.31	68,68,69,69	0
5	DMS	C	8003	4/4	0.91	0.23	58,58,59,61	0
5	DMS	A	8010	4/4	0.91	0.29	58,60,60,61	0
5	DMS	D	8011	4/4	0.92	0.18	40,41,47,51	0
3	NA	D	3102	1/1	0.92	0.08	16,16,16,16	0
5	DMS	C	8017	4/4	0.92	0.15	48,50,50,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	DMS	B	8010	4/4	0.93	0.20	41,45,45,48	0
5	DMS	A	8015	4/4	0.93	0.14	52,54,54,56	0
5	DMS	B	8002	4/4	0.93	0.18	29,32,32,39	0
5	DMS	A	8007	4/4	0.93	0.14	44,46,47,48	0
2	MG	B	3001	1/1	0.93	0.07	13,13,13,13	0
2	MG	D	3001	1/1	0.94	0.09	22,22,22,22	0
5	DMS	B	8011	4/4	0.94	0.24	60,60,60,61	0
5	DMS	C	8012	4/4	0.94	0.24	60,60,62,63	0
5	DMS	B	8015	4/4	0.94	0.16	57,60,60,61	0
3	NA	C	3102	1/1	0.94	0.09	15,15,15,15	0
3	NA	B	3103	1/1	0.94	0.17	28,28,28,28	0
5	DMS	D	8005	4/4	0.94	0.29	42,44,45,47	0
5	DMS	B	8013	4/4	0.94	0.20	46,48,49,49	0
5	DMS	A	8002	4/4	0.94	0.19	30,34,36,37	0
4	IPT	D	2001	15/15	0.94	0.14	17,23,27,30	0
5	DMS	A	8004	4/4	0.94	0.18	48,50,51,52	0
2	MG	C	3001	1/1	0.94	0.09	17,17,17,17	0
5	DMS	D	8009	4/4	0.95	0.16	46,49,50,50	0
3	NA	C	3104	1/1	0.95	0.16	29,29,29,29	0
4	IPT	C	2001	15/15	0.95	0.11	7,14,25,27	0
5	DMS	B	8014	4/4	0.95	0.17	44,48,49,51	0
4	IPT	A	2001	15/15	0.95	0.12	13,18,27,32	0
2	MG	C	3002	1/1	0.95	0.12	13,13,13,13	0
5	DMS	A	8003	4/4	0.95	0.15	29,35,37,38	0
5	DMS	D	8004	4/4	0.95	0.17	40,43,44,47	0
5	DMS	D	8015	4/4	0.95	0.11	45,47,49,54	0
5	DMS	B	8004	4/4	0.95	0.23	27,33,36,37	0
3	NA	B	3102	1/1	0.95	0.10	16,16,16,16	0
5	DMS	B	8001	4/4	0.95	0.21	34,34,35,37	0
5	DMS	B	8009	4/4	0.96	0.18	45,45,46,48	0
5	DMS	C	8002	4/4	0.96	0.14	31,32,33,35	0
5	DMS	C	8016	4/4	0.96	0.16	43,45,48,50	0
3	NA	A	3102	1/1	0.96	0.10	14,14,14,14	0
5	DMS	B	8017	4/4	0.96	0.14	49,50,50,52	0
5	DMS	C	8019	4/4	0.96	0.20	54,56,56,59	0
4	IPT	B	2001	15/15	0.96	0.09	11,15,20,24	0
5	DMS	A	8005	4/4	0.97	0.24	38,40,42,43	0
5	DMS	D	8001	4/4	0.97	0.18	32,34,34,40	0
5	DMS	D	8008	4/4	0.97	0.21	40,47,47,48	0
5	DMS	B	8018	4/4	0.97	0.26	41,43,45,47	0
3	NA	D	3103	1/1	0.97	0.15	29,29,29,29	0
5	DMS	C	8013	4/4	0.97	0.17	65,65,65,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	DMS	D	8003	4/4	0.97	0.14	35,35,36,39	0
5	DMS	A	8009	4/4	0.97	0.17	53,54,55,56	0
5	DMS	C	8009	4/4	0.97	0.15	42,43,44,46	0
3	NA	C	3103	1/1	0.97	0.11	26,26,26,26	0
5	DMS	C	8004	4/4	0.97	0.17	35,38,38,41	0
5	DMS	B	8016	4/4	0.97	0.26	68,68,68,69	0
5	DMS	D	8016	4/4	0.97	0.13	46,47,47,48	0
3	NA	B	3101	1/1	0.97	0.08	15,15,15,15	0
5	DMS	B	8008	4/4	0.97	0.17	35,38,39,42	0
2	MG	A	3001	1/1	0.97	0.05	20,20,20,20	0
5	DMS	A	8008	4/4	0.97	0.21	39,41,43,44	0
5	DMS	C	8008	4/4	0.97	0.16	24,25,27,35	0
2	MG	B	3002	1/1	0.98	0.22	16,16,16,16	0
5	DMS	A	8016	4/4	0.98	0.11	54,55,55,58	0
3	NA	A	3101	1/1	0.98	0.07	16,16,16,16	0
5	DMS	C	8005	4/4	0.98	0.15	48,48,48,50	0
5	DMS	B	8005	4/4	0.98	0.19	34,38,40,41	0
3	NA	D	3101	1/1	0.98	0.07	22,22,22,22	0
5	DMS	B	8003	4/4	0.98	0.18	31,31,32,36	0
3	NA	C	3101	1/1	0.98	0.08	13,13,13,13	0
5	DMS	A	8001	4/4	0.99	0.13	24,26,28,32	0

## 6.5 Other polymers

There are no such residues in this entry.