



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:12 PM GMT

PDB ID : 3WF2  
Title : Crystal structure of human beta-galactosidase in complex with NBT-DGJ  
Authors : Suzuki, H.; Ohto, U.; Shimizu, T.  
Deposited on : 2013-07-16  
Resolution : 2.30 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

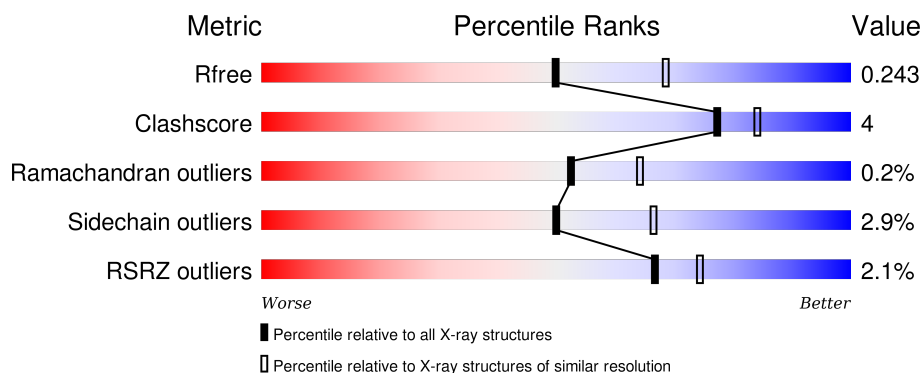
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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	678	<div> <div>2%</div> <div>81%</div> <div>8%</div> <div>11%</div> </div>
1	B	678	<div> <div>%</div> <div>79%</div> <div>9%</div> <div>11%</div> </div>
1	C	678	<div> <div>2%</div> <div>79%</div> <div>10%</div> <div>11%</div> </div>
1	D	678	<div> <div>3%</div> <div>78%</div> <div>10%</div> <div>11%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	702	-	-	-	X
2	NAG	A	703	-	-	-	X
2	NAG	B	703	-	-	-	X
4	NDJ	A	706	-	-	-	X
4	NDJ	B	706	-	-	-	X
6	EDO	A	710	-	-	X	X
6	EDO	B	709	-	-	-	X
6	EDO	C	710	-	-	-	X
6	EDO	D	706	-	-	-	X

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 20707 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	605	Total	C	N	O	S	0	2	0
			4820	3130	794	878	18			
1	B	605	Total	C	N	O	S	0	1	0
			4818	3129	794	878	17			
1	C	603	Total	C	N	O	S	0	4	0
			4805	3121	789	878	17			
1	D	603	Total	C	N	O	S	0	4	0
			4800	3115	789	879	17			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLU	-	EXPRESSION TAG	UNP P16278
A	1	ALA	-	EXPRESSION TAG	UNP P16278
A	2	GLU	-	EXPRESSION TAG	UNP P16278
A	3	ALA	-	EXPRESSION TAG	UNP P16278
A	4	TYR	-	EXPRESSION TAG	UNP P16278
A	5	VAL	-	EXPRESSION TAG	UNP P16278
A	6	GLU	-	EXPRESSION TAG	UNP P16278
A	7	PHE	-	EXPRESSION TAG	UNP P16278
A	8	HIS	-	EXPRESSION TAG	UNP P16278
A	9	HIS	-	EXPRESSION TAG	UNP P16278
A	10	HIS	-	EXPRESSION TAG	UNP P16278
A	11	HIS	-	EXPRESSION TAG	UNP P16278
A	12	HIS	-	EXPRESSION TAG	UNP P16278
A	13	HIS	-	EXPRESSION TAG	UNP P16278
A	14	ASP	-	EXPRESSION TAG	UNP P16278
A	15	TYR	-	EXPRESSION TAG	UNP P16278
A	16	LYS	-	EXPRESSION TAG	UNP P16278
A	17	ASP	-	EXPRESSION TAG	UNP P16278
A	18	ASP	-	EXPRESSION TAG	UNP P16278
A	19	ASP	-	EXPRESSION TAG	UNP P16278
A	20	ASP	-	EXPRESSION TAG	UNP P16278

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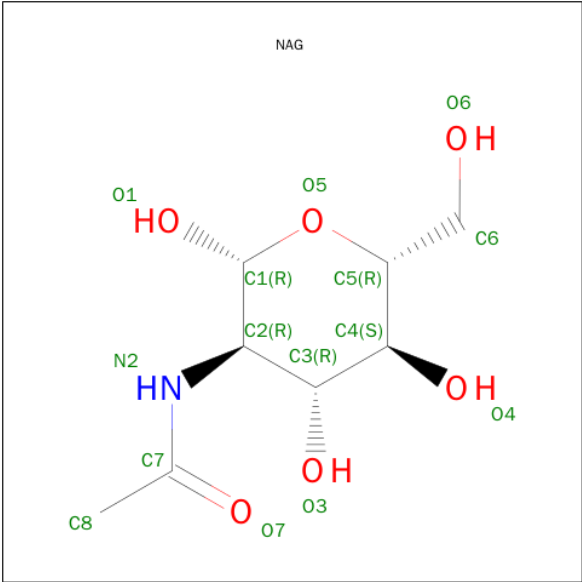
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	LYS	-	EXPRESSION TAG	UNP P16278
A	22	THR	-	EXPRESSION TAG	UNP P16278
A	23	SER	-	EXPRESSION TAG	UNP P16278
B	0	GLU	-	EXPRESSION TAG	UNP P16278
B	1	ALA	-	EXPRESSION TAG	UNP P16278
B	2	GLU	-	EXPRESSION TAG	UNP P16278
B	3	ALA	-	EXPRESSION TAG	UNP P16278
B	4	TYR	-	EXPRESSION TAG	UNP P16278
B	5	VAL	-	EXPRESSION TAG	UNP P16278
B	6	GLU	-	EXPRESSION TAG	UNP P16278
B	7	PHE	-	EXPRESSION TAG	UNP P16278
B	8	HIS	-	EXPRESSION TAG	UNP P16278
B	9	HIS	-	EXPRESSION TAG	UNP P16278
B	10	HIS	-	EXPRESSION TAG	UNP P16278
B	11	HIS	-	EXPRESSION TAG	UNP P16278
B	12	HIS	-	EXPRESSION TAG	UNP P16278
B	13	HIS	-	EXPRESSION TAG	UNP P16278
B	14	ASP	-	EXPRESSION TAG	UNP P16278
B	15	TYR	-	EXPRESSION TAG	UNP P16278
B	16	LYS	-	EXPRESSION TAG	UNP P16278
B	17	ASP	-	EXPRESSION TAG	UNP P16278
B	18	ASP	-	EXPRESSION TAG	UNP P16278
B	19	ASP	-	EXPRESSION TAG	UNP P16278
B	20	ASP	-	EXPRESSION TAG	UNP P16278
B	21	LYS	-	EXPRESSION TAG	UNP P16278
B	22	THR	-	EXPRESSION TAG	UNP P16278
B	23	SER	-	EXPRESSION TAG	UNP P16278
C	0	GLU	-	EXPRESSION TAG	UNP P16278
C	1	ALA	-	EXPRESSION TAG	UNP P16278
C	2	GLU	-	EXPRESSION TAG	UNP P16278
C	3	ALA	-	EXPRESSION TAG	UNP P16278
C	4	TYR	-	EXPRESSION TAG	UNP P16278
C	5	VAL	-	EXPRESSION TAG	UNP P16278
C	6	GLU	-	EXPRESSION TAG	UNP P16278
C	7	PHE	-	EXPRESSION TAG	UNP P16278
C	8	HIS	-	EXPRESSION TAG	UNP P16278
C	9	HIS	-	EXPRESSION TAG	UNP P16278
C	10	HIS	-	EXPRESSION TAG	UNP P16278
C	11	HIS	-	EXPRESSION TAG	UNP P16278
C	12	HIS	-	EXPRESSION TAG	UNP P16278
C	13	HIS	-	EXPRESSION TAG	UNP P16278
C	14	ASP	-	EXPRESSION TAG	UNP P16278

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Chain	Residue	Modelled	Actual	Comment	Reference
C	15	TYR	-	EXPRESSION TAG	UNP P16278
C	16	LYS	-	EXPRESSION TAG	UNP P16278
C	17	ASP	-	EXPRESSION TAG	UNP P16278
C	18	ASP	-	EXPRESSION TAG	UNP P16278
C	19	ASP	-	EXPRESSION TAG	UNP P16278
C	20	ASP	-	EXPRESSION TAG	UNP P16278
C	21	LYS	-	EXPRESSION TAG	UNP P16278
C	22	THR	-	EXPRESSION TAG	UNP P16278
C	23	SER	-	EXPRESSION TAG	UNP P16278
D	0	GLU	-	EXPRESSION TAG	UNP P16278
D	1	ALA	-	EXPRESSION TAG	UNP P16278
D	2	GLU	-	EXPRESSION TAG	UNP P16278
D	3	ALA	-	EXPRESSION TAG	UNP P16278
D	4	TYR	-	EXPRESSION TAG	UNP P16278
D	5	VAL	-	EXPRESSION TAG	UNP P16278
D	6	GLU	-	EXPRESSION TAG	UNP P16278
D	7	PHE	-	EXPRESSION TAG	UNP P16278
D	8	HIS	-	EXPRESSION TAG	UNP P16278
D	9	HIS	-	EXPRESSION TAG	UNP P16278
D	10	HIS	-	EXPRESSION TAG	UNP P16278
D	11	HIS	-	EXPRESSION TAG	UNP P16278
D	12	HIS	-	EXPRESSION TAG	UNP P16278
D	13	HIS	-	EXPRESSION TAG	UNP P16278
D	14	ASP	-	EXPRESSION TAG	UNP P16278
D	15	TYR	-	EXPRESSION TAG	UNP P16278
D	16	LYS	-	EXPRESSION TAG	UNP P16278
D	17	ASP	-	EXPRESSION TAG	UNP P16278
D	18	ASP	-	EXPRESSION TAG	UNP P16278
D	19	ASP	-	EXPRESSION TAG	UNP P16278
D	20	ASP	-	EXPRESSION TAG	UNP P16278
D	21	LYS	-	EXPRESSION TAG	UNP P16278
D	22	THR	-	EXPRESSION TAG	UNP P16278
D	23	SER	-	EXPRESSION TAG	UNP P16278

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
			Total	C	N	O		
			14	8	1	5		
			Total	C	N	O		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
			Total	C	N	O		
			14	8	1	5		
			Total	C	N	O		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
			Total	C	N	O		
			14	8	1	5		
			Total	C	N	O		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
			Total	C	N	O		
			14	8	1	5		
			Total	C	N	O		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
			Total	C	N	O		
			14	8	1	5		
			Total	C	N	O		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
			Total	C	N	O		
			14	8	1	5		
			Total	C	N	O		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
			Total	C	N	O		
			14	8	1	5		
			Total	C	N	O		

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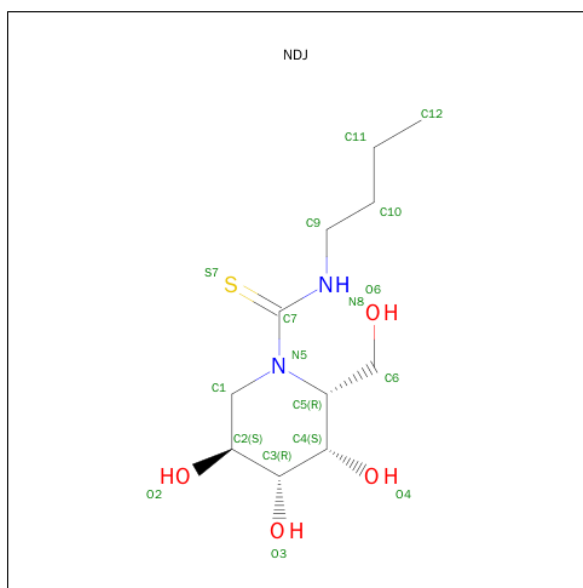
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	A	1	Total	Cl	0	0
			1	1		
3	D	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		

- Molecule 4 is (2R,3S,4R,5S)-N-BUTYL-3,4,5-TRIHYDROXY-2-(HYDROXYMETHYL)PIPERIDINE-1-CARBOTHIOAMIDE (three-letter code: NDJ) (formula: C<sub>11</sub>H<sub>22</sub>N<sub>2</sub>O<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			18	11	2	4	1		
4	B	1	Total	C	N	O	S	0	0
			18	11	2	4	1		
4	C	1	Total	C	N	O	S	0	0
			18	11	2	4	1		

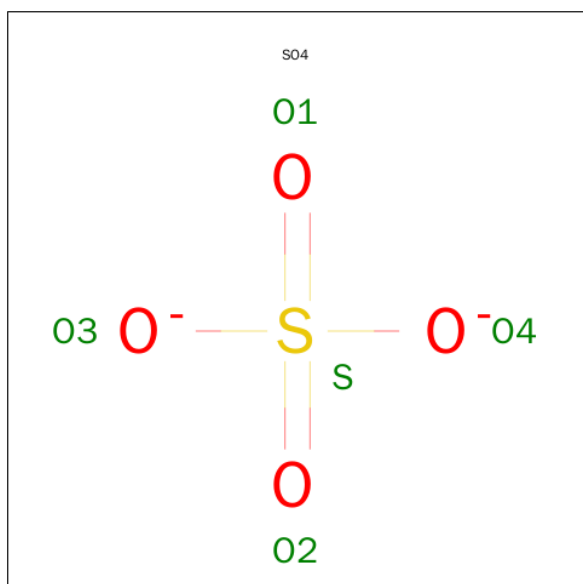
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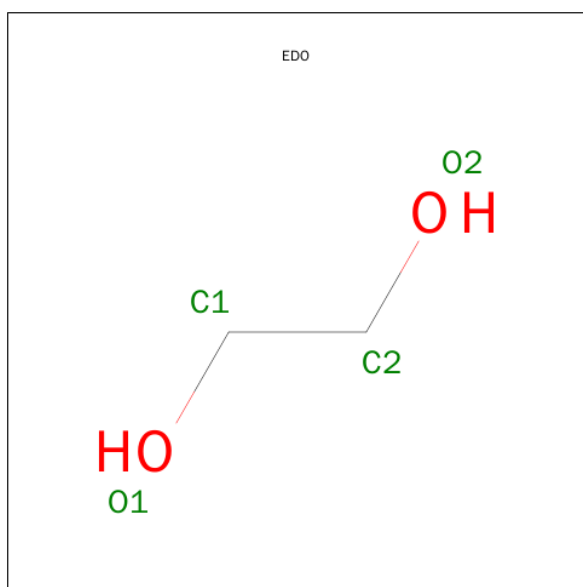
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	D	1	Total	C	N	O	S	0	0
			18	11	2	4	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

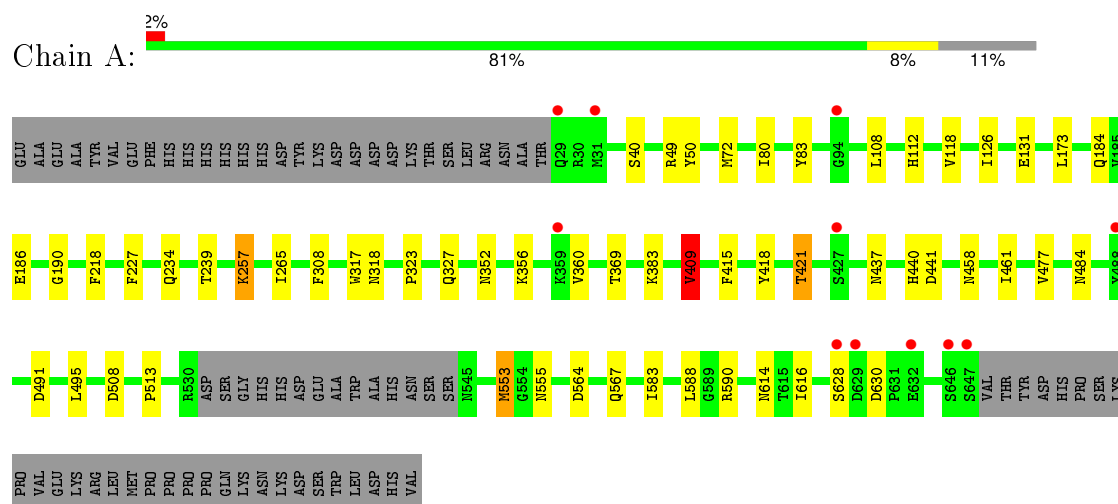
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	288	Total	O	0	0
			288	288		
7	B	288	Total	O	0	0
			288	288		
7	C	279	Total	O	0	0
			279	279		
7	D	237	Total	O	0	0
			237	237		

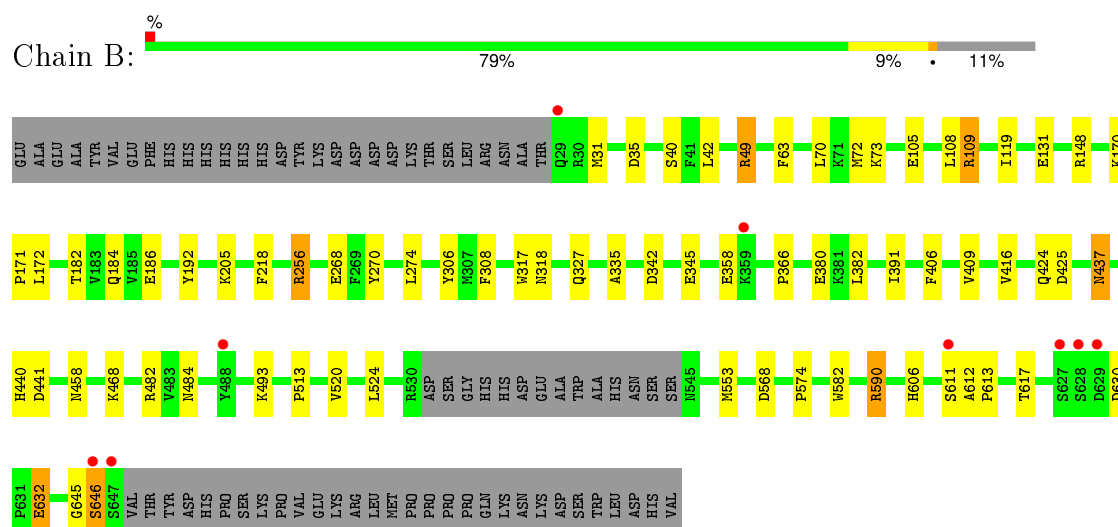
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

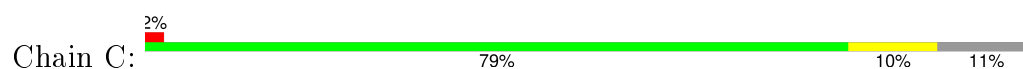
- Molecule 1: Beta-galactosidase

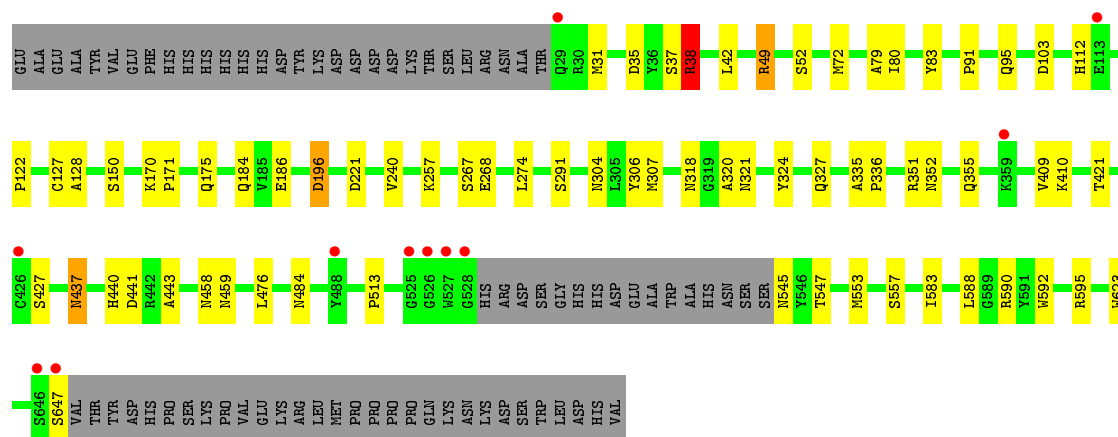


- Molecule 1: Beta-galactosidase

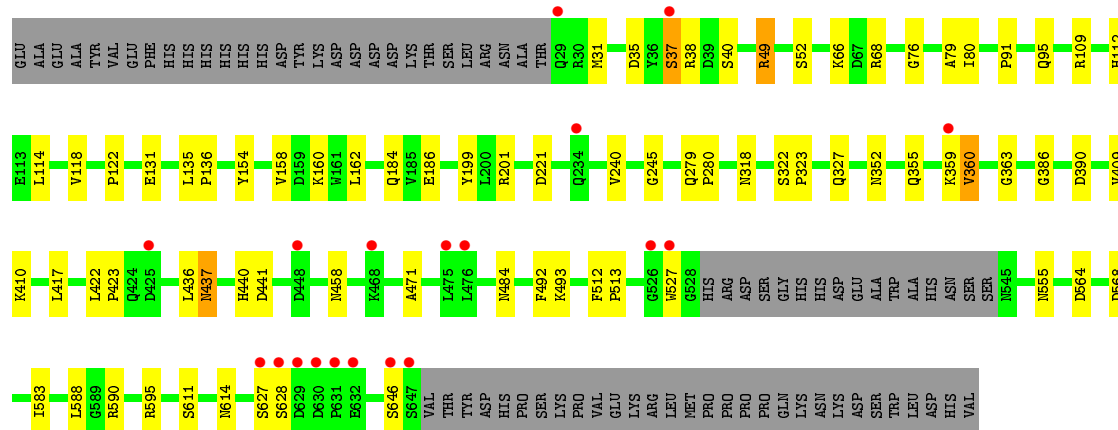
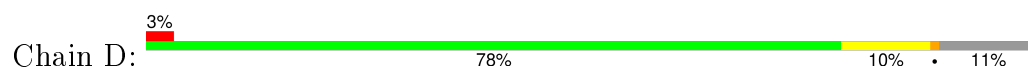


- Molecule 1: Beta-galactosidase





• Molecule 1: Beta-galactosidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.12Å 116.26Å 141.77Å 90.00° 92.41° 90.00°	Depositor
Resolution (Å)	25.98 – 2.30 25.98 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.4 (25.98-2.30) 97.5 (25.98-2.30)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.79 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.175 , 0.240 0.182 , 0.243	Depositor DCC
$R_{free}$ test set	6716 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	29.8	Xtriage
Anisotropy	0.102	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 35.1	EDS
Estimated twinning fraction	0.044 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 133408 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	20707	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 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: SO4, NDJ, EDO, NAG, CL

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	0.80	0/4982	0.87	4/6796 (0.1%)
1	B	0.84	1/4974 (0.0%)	0.90	6/6786 (0.1%)
1	C	0.80	0/4977	0.86	5/6790 (0.1%)
1	D	0.78	0/4971	0.87	6/6782 (0.1%)
All	All	0.81	1/19904 (0.0%)	0.87	21/27154 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	317	TRP	CZ3-CH2	5.12	1.48	1.40

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	38	ARG	NE-CZ-NH1	7.51	124.05	120.30
1	B	342	ASP	CB-CG-OD1	7.22	124.80	118.30
1	A	491	ASP	CB-CG-OD1	6.82	124.44	118.30
1	D	109	ARG	NE-CZ-NH1	6.66	123.63	120.30
1	D	568	ASP	CB-CG-OD1	6.63	124.27	118.30
1	A	491	ASP	CB-CG-OD2	-6.39	112.55	118.30
1	D	68	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	B	109	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	B	49	ARG	CG-CD-NE	-6.09	99.02	111.80
1	B	590	ARG	NE-CZ-NH1	5.83	123.22	120.30
1	A	409	VAL	CB-CA-C	5.79	122.40	111.40
1	D	49	ARG	CG-CD-NE	-5.75	99.73	111.80
1	D	201	ARG	NE-CZ-NH1	5.51	123.06	120.30
1	B	568	ASP	CB-CG-OD2	-5.49	113.36	118.30
1	B	256	ARG	NE-CZ-NH1	-5.37	117.61	120.30
1	A	383	LYS	CB-CA-C	-5.32	99.77	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	196	ASP	CB-CG-OD2	-5.25	113.58	118.30
1	D	595	ARG	NE-CZ-NH1	5.11	122.86	120.30
1	C	49	ARG	CG-CD-NE	-5.07	101.16	111.80
1	C	103	ASP	CB-CG-OD2	-5.05	113.76	118.30
1	C	351	ARG	NE-CZ-NH1	-5.01	117.80	120.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

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	4820	0	4688	34	0
1	B	4818	0	4687	35	0
1	C	4805	0	4670	36	0
1	D	4800	0	4666	37	0
2	A	56	0	52	0	0
2	B	56	0	52	0	0
2	C	56	0	52	0	0
2	D	56	0	52	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	18	0	22	0	0
4	B	18	0	22	1	0
4	C	18	0	22	1	0
4	D	18	0	22	1	0
5	A	10	0	0	1	0
5	B	10	0	0	0	0
5	C	10	0	0	0	0
5	D	10	0	0	0	0
6	A	8	0	12	4	0
6	B	8	0	12	0	0
6	C	8	0	12	0	0
6	D	8	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	288	0	0	5	0
7	B	288	0	0	5	0
7	C	279	0	0	3	0
7	D	237	0	0	4	0
All	All	20707	0	19055	142	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (142) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:VAL:HG13	1:A:513:PRO:HG3	1.62	0.81
1:B:437:ASN:HD21	1:B:458:ASN:H	1.29	0.80
1:D:555:ASN:HB2	2:D:704:NAG:O5	1.83	0.78
1:D:35:ASP:HB3	7:D:1006:HOH:O	1.82	0.78
1:C:318:ASN:HD21	1:C:590:ARG:HH21	1.31	0.77
1:C:437:ASN:HD21	1:C:458:ASN:H	1.33	0.74
1:A:437:ASN:HD21	1:A:458:ASN:H	1.37	0.72
1:B:318:ASN:HD21	1:B:590:ARG:HH21	1.39	0.71
1:D:437:ASN:HD21	1:D:458:ASN:H	1.36	0.70
1:C:112:HIS:CD2	7:C:1050:HOH:O	2.48	0.67
1:B:409:VAL:HG13	1:B:513:PRO:HG3	1.75	0.66
1:C:38:ARG:HH11	1:C:38:ARG:HG3	1.60	0.66
1:B:553:MET:HE3	1:B:617:THR:HG23	1.78	0.66
1:A:327:GLN:HE22	1:A:484:ASN:HD21	1.47	0.62
1:B:612:ALA:HB1	1:B:613:PRO:HD2	1.82	0.61
1:D:91:PRO:HD2	1:D:95:GLN:O	2.01	0.61
1:A:257:LYS:HD3	7:A:1059:HOH:O	2.01	0.61
1:B:382:LEU:HD23	1:B:524:LEU:HD12	1.81	0.61
1:A:317:TRP:CZ2	6:A:710:EDO:H11	2.37	0.59
1:A:112:HIS:HD2	7:A:1063:HOH:O	1.84	0.59
1:D:327:GLN:HE22	1:D:484:ASN:HD21	1.50	0.59
1:B:574:PRO:CD	7:B:1068:HOH:O	2.52	0.58
1:B:574:PRO:HD2	7:B:1068:HOH:O	2.03	0.57
1:B:105:GLU:O	1:B:109:ARG:HG3	2.04	0.57
1:C:583:ILE:HD12	1:C:588:LEU:HD11	1.87	0.56
1:D:422:LEU:HD12	1:D:471:ALA:HB3	1.87	0.56
1:A:421:THR:HB	7:A:1070:HOH:O	2.06	0.55
1:D:112:HIS:HD2	7:D:1005:HOH:O	1.89	0.55
1:D:352[A]:ASN:OD1	1:D:355:GLN:NE2	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:318:ASN:HD21	1:B:590:ARG:NH2	2.03	0.55
1:C:112:HIS:HD2	7:C:1050:HOH:O	1.87	0.55
1:C:409:VAL:HG13	1:C:513:PRO:HG3	1.88	0.54
1:B:73:LYS:HE3	7:B:1059:HOH:O	2.07	0.54
1:A:628:SER:O	1:A:630:ASP:N	2.40	0.54
1:C:91:PRO:HD2	1:C:95:GLN:O	2.08	0.53
1:D:555:ASN:CB	2:D:704:NAG:O5	2.54	0.53
1:C:122:PRO:HD2	1:C:184:GLN:O	2.08	0.53
1:A:108:LEU:HD22	1:A:118:VAL:HG11	1.90	0.53
1:B:612:ALA:HB1	1:B:613:PRO:CD	2.39	0.53
1:C:268:GLU:OE1	4:C:705:NDJ:H5	2.09	0.53
1:C:327:GLN:HE22	1:C:484:ASN:HD21	1.55	0.52
1:C:83:TYR:CE2	1:C:128:ALA:HB2	2.44	0.52
1:C:170:LYS:HE3	7:C:1061:HOH:O	2.09	0.51
1:D:318:ASN:HD21	1:D:590:ARG:HH21	1.57	0.51
1:D:417:LEU:HB3	1:D:512:PHE:HB2	1.92	0.51
1:A:369:THR:HG22	1:D:564:ASP:OD2	2.10	0.51
1:D:52:SER:HA	1:D:79:ALA:O	2.11	0.50
1:D:38:ARG:HB2	7:D:1006:HOH:O	2.12	0.50
1:C:267:SER:O	1:C:304:ASN:HB3	2.12	0.50
1:A:72:MET:HG2	1:A:308:PHE:CD2	2.48	0.49
1:C:150:SER:OG	1:C:196:ASP:OD2	2.30	0.49
1:A:83:TYR:CD1	1:A:126:ILE:HB	2.48	0.48
1:D:555:ASN:HB2	2:D:704:NAG:H61	1.95	0.48
1:A:49:ARG:HG2	1:A:50:TYR:N	2.18	0.48
1:A:323:PRO:HA	5:A:708:SO4:O3	2.13	0.48
1:B:630:ASP:OD1	1:B:632:GLU:HB3	2.13	0.48
1:B:406:PHE:CZ	1:B:416:VAL:HG23	2.49	0.48
1:C:170:LYS:HB3	1:C:171:PRO:CD	2.43	0.47
1:B:553:MET:HE3	1:B:617:THR:CG2	2.43	0.47
1:B:268:GLU:OE1	4:B:706:NDJ:H5	2.15	0.47
1:D:66:LYS:HE3	1:D:114:LEU:HD21	1.97	0.47
1:B:308:PHE:O	1:B:335:ALA:HB1	2.15	0.47
1:A:553:MET:HA	1:A:616:ILE:O	2.14	0.47
1:A:564:ASP:O	1:A:567:GLN:HG3	2.15	0.47
1:A:184:GLN:HB2	1:A:218:PHE:CZ	2.51	0.46
1:B:184:GLN:HB2	1:B:218:PHE:CZ	2.51	0.46
1:B:482:ARG:HG2	1:B:493:LYS:HE2	1.98	0.46
1:B:119:ILE:HG23	1:B:182:THR:OG1	2.15	0.46
1:A:317:TRP:CE2	6:A:710:EDO:C1	2.98	0.46
1:D:440:HIS:HA	1:D:441:ASP:HA	1.77	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:ASN:HB3	7:A:1076:HOH:O	2.16	0.46
1:B:35:ASP:HB2	1:B:42:LEU:HG	1.97	0.45
1:D:363:GLY:HA3	7:D:1000:HOH:O	2.15	0.45
1:C:592:TRP:CE2	1:C:595:ARG:HG3	2.52	0.45
1:B:327:GLN:HE22	1:B:484:ASN:HD21	1.64	0.45
1:A:190:GLY:HA3	1:A:227:PHE:O	2.16	0.45
1:C:38:ARG:CG	1:C:38:ARG:HH11	2.27	0.45
1:C:35:ASP:OD2	1:C:38:ARG:HB2	2.17	0.45
1:A:440:HIS:HA	1:A:441:ASP:HA	1.80	0.45
1:C:592:TRP:CZ2	1:C:595:ARG:HG3	2.52	0.44
1:B:318:ASN:ND2	1:B:590:ARG:HH21	2.10	0.44
1:D:279:GLN:HE21	1:D:280:PRO:HD2	1.83	0.44
1:D:583:ILE:HD12	1:D:588:LEU:HD11	2.00	0.44
1:D:322:SER:HA	1:D:323:PRO:C	2.38	0.44
1:A:80:ILE:HD11	1:A:118:VAL:HG22	1.99	0.44
1:D:492:PHE:O	1:D:493:LYS:HB2	2.17	0.44
1:D:154:TYR:HE2	1:D:199:TYR:CE1	2.35	0.44
1:D:76:GLY:CA	1:D:360:VAL:HG22	2.48	0.44
1:A:239:THR:HG22	1:A:265:ILE:HD12	2.00	0.44
1:C:72:MET:HE2	1:C:80:ILE:HG22	2.00	0.44
1:A:555:ASN:HA	1:A:614:ASN:O	2.19	0.43
1:B:170:LYS:N	1:B:171:PRO:HD2	2.34	0.43
1:D:158:VAL:O	1:D:162:LEU:HG	2.18	0.43
1:C:52:SER:HA	1:C:79:ALA:O	2.18	0.43
1:D:555:ASN:HA	1:D:614:ASN:O	2.18	0.43
1:D:80:ILE:HD11	1:D:118:VAL:HG22	1.99	0.43
1:A:437:ASN:ND2	1:A:458:ASN:H	2.12	0.42
4:D:708:NDJ:N8	4:D:708:NDJ:H9	2.34	0.42
1:C:320:ALA:H	1:C:484:ASN:HD22	1.67	0.42
1:B:63:PHE:CD2	1:B:606:HIS:HB3	2.55	0.42
1:C:409:VAL:O	1:C:410:LYS:HB2	2.20	0.42
1:B:645:GLY:O	1:B:646:SER:C	2.57	0.42
1:B:72:MET:HG2	1:B:308:PHE:CD2	2.55	0.42
1:A:318:ASN:HD21	1:A:590:ARG:HH21	1.67	0.42
1:A:415:PHE:HA	1:A:477:VAL:O	2.20	0.42
1:C:547:THR:OG1	1:C:623:TRP:HA	2.20	0.42
1:B:108:LEU:HD12	1:B:172:LEU:HD13	2.01	0.42
1:D:135:LEU:HA	1:D:136:PRO:HD3	1.91	0.41
1:B:31:MET:HA	7:B:890:HOH:O	2.20	0.41
1:A:583:ILE:HD12	1:A:588:LEU:HD11	2.02	0.41
1:B:148:ARG:HB2	1:B:192:TYR:CE2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:35:ASP:HB2	1:C:42:LEU:HG	2.02	0.41
1:A:173:LEU:HA	1:A:173:LEU:HD23	1.86	0.41
1:D:122:PRO:HD2	1:D:184:GLN:O	2.20	0.41
1:D:245:GLY:H	6:D:706:EDO:H22	1.84	0.41
1:C:352[A]:ASN:OD1	1:C:355:GLN:NE2	2.54	0.41
1:D:221:ASP:O	1:D:240:VAL:HA	2.21	0.41
1:C:440:HIS:HA	1:C:441:ASP:HA	1.85	0.41
1:D:437:ASN:OD1	1:D:458:ASN:OD1	2.39	0.41
1:A:317:TRP:CE2	6:A:710:EDO:H11	2.56	0.41
1:C:409:VAL:O	1:C:410:LYS:CB	2.69	0.41
1:B:70:LEU:HD23	1:B:366:PRO:HG2	2.02	0.41
1:C:321:ASN:O	1:C:324:TYR:HA	2.21	0.41
1:A:317:TRP:CZ2	6:A:710:EDO:C1	3.02	0.41
1:B:520:VAL:HG11	1:B:582:TRP:CG	2.56	0.41
1:C:221:ASP:O	1:C:240:VAL:HA	2.20	0.41
1:C:306:TYR:HA	1:C:307:MET:HA	1.93	0.41
1:A:461:ILE:HG21	1:A:461:ILE:HD13	1.87	0.41
1:B:270:TYR:HA	1:B:306:TYR:O	2.21	0.41
1:B:440:HIS:HA	1:B:441:ASP:HA	1.86	0.41
1:A:112:HIS:CD2	7:A:1063:HOH:O	2.65	0.40
1:C:409:VAL:CG1	1:C:409:VAL:O	2.68	0.40
1:D:37:SER:HB2	1:D:38:ARG:HG2	2.02	0.40
1:D:409:VAL:O	1:D:410:LYS:HB2	2.20	0.40
1:C:335:ALA:HB1	1:C:336:PRO:HD2	2.03	0.40
1:A:418:TYR:CD2	1:A:495:LEU:HD12	2.55	0.40
1:D:436:LEU:O	1:D:437:ASN:HB3	2.21	0.40
1:B:574:PRO:HG2	7:B:1068:HOH:O	2.22	0.40
1:D:409:VAL:HG13	1:D:513:PRO:HG3	2.03	0.40
1:C:443:ALA:HA	1:C:476:LEU:O	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	603/678 (89%)	570 (94%)	32 (5%)	1 (0%)	52	64
1	B	602/678 (89%)	575 (96%)	26 (4%)	1 (0%)	52	64
1	C	603/678 (89%)	576 (96%)	26 (4%)	1 (0%)	52	64
1	D	603/678 (89%)	569 (94%)	32 (5%)	2 (0%)	46	57
All	All	2411/2712 (89%)	2290 (95%)	116 (5%)	5 (0%)	52	64

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	646	SER
1	A	508	ASP
1	C	127	CYS
1	D	423	PRO
1	D	386	GLY

### 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	520/585 (89%)	510 (98%)	10 (2%)	65	81
1	B	519/585 (89%)	502 (97%)	17 (3%)	45	61
1	C	520/585 (89%)	502 (96%)	18 (4%)	43	58
1	D	520/585 (89%)	503 (97%)	17 (3%)	45	61
All	All	2079/2340 (89%)	2017 (97%)	62 (3%)	50	65

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	SER
1	A	131	GLU
1	A	186	GLU
1	A	234	GLN
1	A	257	LYS
1	A	356	LYS

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Mol	Chain	Res	Type
1	A	360	VAL
1	A	409	VAL
1	A	421	THR
1	A	553	MET
1	B	40	SER
1	B	49	ARG
1	B	131	GLU
1	B	186	GLU
1	B	205	LYS
1	B	256	ARG
1	B	274	LEU
1	B	345	GLU
1	B	358	GLU
1	B	380	GLU
1	B	391	ILE
1	B	424	GLN
1	B	425	ASP
1	B	437	ASN
1	B	468	LYS
1	B	611	SER
1	B	632	GLU
1	C	31	MET
1	C	37	SER
1	C	38	ARG
1	C	49	ARG
1	C	175[A]	GLN
1	C	175[B]	GLN
1	C	186	GLU
1	C	257	LYS
1	C	274	LEU
1	C	291	SER
1	C	421	THR
1	C	427	SER
1	C	437	ASN
1	C	459	ASN
1	C	545	ASN
1	C	553	MET
1	C	557	SER
1	C	647	SER
1	D	31	MET
1	D	37	SER
1	D	40	SER

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Mol	Chain	Res	Type
1	D	49	ARG
1	D	131	GLU
1	D	160	LYS
1	D	186	GLU
1	D	359	LYS
1	D	360	VAL
1	D	390	ASP
1	D	437	ASN
1	D	527	TRP
1	D	611[A]	SER
1	D	611[B]	SER
1	D	627	SER
1	D	628	SER
1	D	646	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	HIS
1	A	112	HIS
1	A	318	ASN
1	A	412	HIS
1	A	437	ASN
1	A	484	ASN
1	A	529	HIS
1	B	95	GLN
1	B	102	HIS
1	B	318	ASN
1	B	355	GLN
1	B	437	ASN
1	B	484	ASN
1	C	102	HIS
1	C	279	GLN
1	C	318	ASN
1	C	355	GLN
1	C	437	ASN
1	C	484	ASN
1	D	234	GLN
1	D	279	GLN
1	D	318	ASN
1	D	321	ASN
1	D	355	GLN

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Mol	Chain	Res	Type
1	D	437	ASN
1	D	484	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 40 ligands modelled in this entry, 4 are monoatomic - leaving 36 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	$\# Z  > 2$
2	NAG	A	701	1	14,14,15	0.96	2 (14%)	15,19,21	1.34	2 (13%)
2	NAG	A	702	1	14,14,15	0.48	0	15,19,21	2.47	3 (20%)
2	NAG	A	703	1	14,14,15	1.45	3 (21%)	15,19,21	3.63	8 (53%)
2	NAG	A	704	1	14,14,15	0.81	0	15,19,21	1.60	3 (20%)
4	NDJ	A	706	-	17,18,18	1.40	5 (29%)	18,24,24	1.35	3 (16%)
5	SO4	A	707	-	4,4,4	0.63	0	6,6,6	0.17	0
5	SO4	A	708	-	4,4,4	0.91	0	6,6,6	0.88	0
6	EDO	A	709	-	3,3,3	0.44	0	2,2,2	0.80	0
6	EDO	A	710	-	3,3,3	0.38	0	2,2,2	0.22	0
2	NAG	B	701	1	14,14,15	0.58	0	15,19,21	1.43	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	B	702	1	14,14,15	0.79	0	15,19,21	1.50	2 (13%)
2	NAG	B	703	1	14,14,15	0.85	1 (7%)	15,19,21	2.50	3 (20%)
2	NAG	B	704	1	14,14,15	0.48	0	15,19,21	0.73	0
4	NDJ	B	706	-	17,18,18	1.78	4 (23%)	18,24,24	1.59	4 (22%)
5	SO4	B	707	-	4,4,4	0.52	0	6,6,6	0.23	0
5	SO4	B	708	-	4,4,4	0.72	0	6,6,6	1.11	0
6	EDO	B	709	-	3,3,3	0.43	0	2,2,2	0.68	0
6	EDO	B	710	-	3,3,3	0.51	0	2,2,2	0.29	0
2	NAG	C	701	1	14,14,15	0.65	0	15,19,21	1.97	3 (20%)
2	NAG	C	702	1	14,14,15	0.71	1 (7%)	15,19,21	1.87	2 (13%)
2	NAG	C	703	1	14,14,15	0.75	0	15,19,21	1.64	4 (26%)
2	NAG	C	704	1	14,14,15	0.73	0	15,19,21	1.55	5 (33%)
4	NDJ	C	705	-	17,18,18	1.26	1 (5%)	18,24,24	1.82	4 (22%)
5	SO4	C	706	-	4,4,4	0.51	0	6,6,6	0.25	0
5	SO4	C	707	-	4,4,4	0.76	0	6,6,6	0.70	0
6	EDO	C	709	-	3,3,3	0.75	0	2,2,2	0.27	0
6	EDO	C	710	-	3,3,3	0.61	0	2,2,2	0.31	0
2	NAG	D	701	1	14,14,15	0.28	0	15,19,21	0.54	0
2	NAG	D	703	1	14,14,15	0.63	0	15,19,21	1.68	3 (20%)
2	NAG	D	704	1	14,14,15	0.64	0	15,19,21	2.63	3 (20%)
2	NAG	D	705	1	14,14,15	0.91	0	15,19,21	1.73	5 (33%)
6	EDO	D	706	-	3,3,3	0.30	0	2,2,2	1.80	1 (50%)
6	EDO	D	707	-	3,3,3	0.58	0	2,2,2	0.26	0
4	NDJ	D	708	-	17,18,18	1.28	1 (5%)	18,24,24	1.62	3 (16%)
5	SO4	D	709	-	4,4,4	0.38	0	6,6,6	0.69	0
5	SO4	D	710	-	4,4,4	0.87	0	6,6,6	1.33	1 (16%)

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	NAG	A	701	1	-	0/6/23/26	0/1/1/1
2	NAG	A	702	1	-	0/6/23/26	0/1/1/1
2	NAG	A	703	1	-	0/6/23/26	0/1/1/1
2	NAG	A	704	1	-	0/6/23/26	0/1/1/1
4	NDJ	A	706	-	-	0/10/31/31	0/1/1/1
5	SO4	A	707	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	SO4	A	708	-	-	0/0/0/0	0/0/0/0
6	EDO	A	709	-	-	0/1/1/1	0/0/0/0
6	EDO	A	710	-	-	0/1/1/1	0/0/0/0
2	NAG	B	701	1	-	0/6/23/26	0/1/1/1
2	NAG	B	702	1	-	0/6/23/26	0/1/1/1
2	NAG	B	703	1	-	0/6/23/26	0/1/1/1
2	NAG	B	704	1	-	0/6/23/26	0/1/1/1
4	NDJ	B	706	-	-	0/10/31/31	0/1/1/1
5	SO4	B	707	-	-	0/0/0/0	0/0/0/0
5	SO4	B	708	-	-	0/0/0/0	0/0/0/0
6	EDO	B	709	-	-	0/1/1/1	0/0/0/0
6	EDO	B	710	-	-	0/1/1/1	0/0/0/0
2	NAG	C	701	1	-	0/6/23/26	0/1/1/1
2	NAG	C	702	1	-	0/6/23/26	0/1/1/1
2	NAG	C	703	1	-	0/6/23/26	0/1/1/1
2	NAG	C	704	1	-	0/6/23/26	0/1/1/1
4	NDJ	C	705	-	-	0/10/31/31	0/1/1/1
5	SO4	C	706	-	-	0/0/0/0	0/0/0/0
5	SO4	C	707	-	-	0/0/0/0	0/0/0/0
6	EDO	C	709	-	-	0/1/1/1	0/0/0/0
6	EDO	C	710	-	-	0/1/1/1	0/0/0/0
2	NAG	D	701	1	-	0/6/23/26	0/1/1/1
2	NAG	D	703	1	-	0/6/23/26	0/1/1/1
2	NAG	D	704	1	-	0/6/23/26	0/1/1/1
2	NAG	D	705	1	-	0/6/23/26	0/1/1/1
6	EDO	D	706	-	-	0/1/1/1	0/0/0/0
6	EDO	D	707	-	-	0/1/1/1	0/0/0/0
4	NDJ	D	708	-	-	0/10/31/31	0/1/1/1
5	SO4	D	709	-	-	0/0/0/0	0/0/0/0
5	SO4	D	710	-	-	0/0/0/0	0/0/0/0

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	708	NDJ	C7-S7	-4.33	1.60	1.67
4	B	706	NDJ	C7-S7	-4.07	1.61	1.67
4	C	705	NDJ	C7-S7	-3.78	1.61	1.67
4	A	706	NDJ	C7-S7	-2.60	1.63	1.67
2	A	701	NAG	O5-C1	-2.58	1.39	1.43
2	A	701	NAG	C2-N2	-2.08	1.42	1.46
4	A	706	NDJ	C2-C3	2.11	1.55	1.52
4	A	706	NDJ	C7-N8	2.13	1.37	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	703	NAG	C1-C2	2.15	1.55	1.52
2	B	703	NAG	C1-C2	2.17	1.55	1.52
2	C	702	NAG	C1-C2	2.21	1.55	1.52
4	A	706	NDJ	C5-N5	2.39	1.51	1.48
2	A	703	NAG	C3-C2	2.50	1.58	1.52
4	B	706	NDJ	C7-N8	2.51	1.38	1.34
4	A	706	NDJ	C1-C2	2.55	1.56	1.52
4	B	706	NDJ	C1-C2	3.10	1.56	1.52
4	B	706	NDJ	C2-C3	3.36	1.57	1.52
2	A	703	NAG	C2-N2	3.45	1.52	1.46

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	703	NAG	O7-C7-C8	-5.05	112.80	122.06
2	A	703	NAG	O3-C3-C4	-4.70	99.75	110.34
4	C	705	NDJ	C1-N5-C5	-4.23	103.70	115.89
4	C	705	NDJ	S7-C7-N8	-3.77	114.54	122.47
2	D	705	NAG	C3-C4-C5	-3.65	103.83	110.20
4	B	706	NDJ	C1-N5-C5	-3.30	106.38	115.89
4	D	708	NDJ	S7-C7-N8	-3.26	115.61	122.47
4	D	708	NDJ	C1-N5-C5	-3.08	107.01	115.89
2	D	703	NAG	C2-N2-C7	-2.91	119.30	123.04
5	D	710	SO4	O2-S-O1	-2.84	100.48	109.50
2	D	705	NAG	O4-C4-C3	-2.84	103.94	110.34
4	A	706	NDJ	C1-N5-C5	-2.84	107.70	115.89
4	A	706	NDJ	S7-C7-N8	-2.69	116.82	122.47
4	B	706	NDJ	S7-C7-N8	-2.67	116.85	122.47
2	C	704	NAG	O7-C7-C8	-2.58	117.32	122.06
2	D	704	NAG	C4-C3-C2	-2.58	107.22	111.23
2	A	701	NAG	C8-C7-N2	-2.58	111.18	116.11
2	B	703	NAG	C6-C5-C4	-2.53	106.77	113.02
2	D	703	NAG	C3-C4-C5	-2.52	105.80	110.20
2	D	704	NAG	O7-C7-C8	-2.43	117.60	122.06
6	D	706	EDO	O2-C2-C1	-2.41	95.28	112.54
2	A	702	NAG	C6-C5-C4	-2.37	107.16	113.02
2	C	702	NAG	C2-N2-C7	-2.30	120.08	123.04
2	D	705	NAG	C2-N2-C7	-2.28	120.11	123.04
2	C	703	NAG	O3-C3-C2	-2.27	104.61	109.11
2	A	702	NAG	O6-C6-C5	-2.19	104.08	111.33
2	A	703	NAG	C4-C3-C2	-2.18	107.83	111.23
2	B	701	NAG	O6-C6-C5	-2.06	104.53	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	705	NAG	C3-C2-N2	-2.02	105.71	110.56
2	C	704	NAG	O3-C3-C2	-2.02	105.12	109.11
4	C	705	NDJ	C3-C4-C5	2.03	114.12	111.39
2	A	701	NAG	O7-C7-N2	2.04	126.02	121.86
2	B	702	NAG	O7-C7-N2	2.07	126.08	121.86
2	C	704	NAG	O7-C7-N2	2.11	126.16	121.86
2	C	704	NAG	O5-C5-C6	2.31	112.35	107.35
2	C	703	NAG	C8-C7-N2	2.31	120.53	116.11
2	A	704	NAG	C4-C3-C2	2.32	114.83	111.23
2	A	704	NAG	C3-C4-C5	2.38	114.35	110.20
2	C	704	NAG	C1-O5-C5	2.56	115.49	112.25
2	D	705	NAG	C1-O5-C5	2.59	115.53	112.25
4	A	706	NDJ	C1-C2-C3	2.71	113.16	110.21
4	B	706	NDJ	N8-C7-N5	2.76	121.34	115.79
4	B	706	NDJ	O3-C3-C2	2.80	115.06	110.00
2	A	703	NAG	C8-C7-N2	2.89	121.64	116.11
2	C	703	NAG	C1-O5-C5	3.03	116.09	112.25
2	C	703	NAG	O3-C3-C4	3.12	117.36	110.34
4	C	705	NDJ	N8-C7-N5	3.47	122.76	115.79
2	B	703	NAG	C3-C4-C5	3.48	116.27	110.20
2	C	701	NAG	C3-C4-C5	3.76	116.75	110.20
4	D	708	NDJ	N8-C7-N5	3.77	123.36	115.79
2	C	701	NAG	C1-O5-C5	3.82	117.10	112.25
2	B	701	NAG	C1-O5-C5	3.90	117.20	112.25
2	C	701	NAG	C4-C3-C2	4.18	117.72	111.23
2	B	702	NAG	C1-O5-C5	4.22	117.60	112.25
2	A	703	NAG	C2-N2-C7	4.27	128.53	123.04
2	D	703	NAG	C1-O5-C5	4.38	117.81	112.25
2	A	704	NAG	C1-O5-C5	4.47	117.93	112.25
2	A	703	NAG	C3-C2-N2	5.38	123.45	110.56
2	C	702	NAG	C1-O5-C5	5.63	119.39	112.25
2	A	703	NAG	O3-C3-C2	5.78	120.57	109.11
2	A	703	NAG	C1-O5-C5	6.88	120.98	112.25
2	B	703	NAG	C1-O5-C5	7.86	122.22	112.25
2	A	702	NAG	C1-O5-C5	8.39	122.89	112.25
2	D	704	NAG	C1-O5-C5	8.58	123.14	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	708	SO4	1	0
6	A	710	EDO	4	0
4	B	706	NDJ	1	0
4	C	705	NDJ	1	0
2	D	704	NAG	3	0
6	D	706	EDO	1	0
4	D	708	NDJ	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	605/678 (89%)	-0.19	11 (1%) 71 78	18, 29, 48, 104	0
1	B	605/678 (89%)	-0.24	9 (1%) 76 81	16, 27, 46, 93	0
1	C	603/678 (88%)	-0.12	11 (1%) 71 78	17, 29, 56, 87	0
1	D	603/678 (88%)	-0.08	19 (3%) 51 60	17, 30, 58, 97	0
All	All	2416/2712 (89%)	-0.16	50 (2%) 67 74	16, 29, 54, 104	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	647	SER	6.6
1	C	527	TRP	6.4
1	A	647	SER	5.6
1	A	29	GLN	5.5
1	D	629	ASP	5.1
1	C	525	GLY	5.0
1	B	627	SER	5.0
1	D	628	SER	4.9
1	A	629	ASP	4.9
1	B	646	SER	4.8
1	D	29	GLN	4.8
1	B	29	GLN	4.4
1	D	527	TRP	4.2
1	A	646	SER	4.1
1	D	647	SER	4.0
1	C	528	GLY	4.0
1	D	526	GLY	3.7
1	B	629	ASP	3.7
1	B	628	SER	3.6
1	C	29	GLN	3.5
1	A	628	SER	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	359	LYS	3.1
1	D	646	SER	3.0
1	C	426	CYS	2.9
1	D	37	SER	2.9
1	C	526	GLY	2.9
1	D	632	GLU	2.8
1	C	488	TYR	2.8
1	C	359	LYS	2.8
1	D	425	ASP	2.7
1	B	611	SER	2.7
1	D	476	LEU	2.5
1	A	632	GLU	2.5
1	B	488	TYR	2.5
1	D	448	ASP	2.5
1	D	234	GLN	2.5
1	D	630	ASP	2.4
1	A	359	LYS	2.3
1	D	627	SER	2.3
1	D	468	LYS	2.3
1	A	31	MET	2.2
1	A	94	GLY	2.2
1	C	113	GLU	2.2
1	D	631	PRO	2.1
1	C	647	SER	2.1
1	B	359	LYS	2.1
1	C	646	SER	2.1
1	D	475	LEU	2.1
1	A	488	TYR	2.1
1	A	427	SER	2.1

## 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
6	EDO	A	710	4/4	0.96	0.19	5.29	34,45,46,47	0
6	EDO	D	706	4/4	0.93	0.15	4.43	35,38,39,40	0
2	NAG	A	703	14/15	0.83	0.18	4.28	43,50,55,55	0
6	EDO	C	710	4/4	0.88	0.22	4.09	37,46,54,55	0
2	NAG	A	702	14/15	0.88	0.19	3.44	25,36,42,45	0
4	NDJ	B	706	18/18	0.88	0.19	3.04	22,29,81,83	0
6	EDO	B	709	4/4	0.77	0.21	2.87	47,51,53,55	0
4	NDJ	A	706	18/18	0.85	0.20	2.52	25,29,59,66	0
2	NAG	B	703	14/15	0.90	0.18	2.36	46,52,61,61	0
5	SO4	D	710	5/5	0.88	0.20	1.93	54,59,70,70	0
2	NAG	C	703	14/15	0.88	0.23	1.54	40,52,55,58	0
6	EDO	A	709	4/4	0.88	0.17	1.49	42,47,53,58	0
5	SO4	A	708	5/5	0.94	0.19	1.38	47,48,59,62	0
4	NDJ	D	708	18/18	0.85	0.18	1.36	27,35,58,77	0
2	NAG	B	702	14/15	0.93	0.14	1.30	31,35,38,39	0
2	NAG	D	703	14/15	0.87	0.22	1.10	42,53,62,63	0
5	SO4	B	708	5/5	0.93	0.19	1.09	57,58,61,66	0
2	NAG	C	702	14/15	0.87	0.22	1.01	53,58,61,63	0
4	NDJ	C	705	18/18	0.88	0.16	0.94	19,31,63,77	0
2	NAG	D	705	14/15	0.94	0.14	0.91	32,36,39,44	0
5	SO4	C	707	5/5	0.93	0.18	0.60	58,58,62,62	0
2	NAG	C	704	14/15	0.94	0.10	-0.03	27,31,34,34	0
2	NAG	B	701	14/15	0.96	0.10	-0.57	35,40,49,49	0
2	NAG	A	701	14/15	0.95	0.10	-0.61	26,30,34,38	0
3	CL	D	702	1/1	0.99	0.12	-1.21	22,22,22,22	0
3	CL	C	708	1/1	1.00	0.10	-2.31	19,19,19,19	0
3	CL	A	705	1/1	0.99	0.05	-2.80	21,21,21,21	0
3	CL	B	705	1/1	0.99	0.09	-3.55	21,21,21,21	0
6	EDO	B	710	4/4	0.90	0.19	-	43,43,47,48	0
2	NAG	D	704	14/15	0.58	0.34	-	72,93,100,101	0
6	EDO	D	707	4/4	0.95	0.16	-	35,44,47,49	0
5	SO4	C	706	5/5	0.93	0.42	-	70,73,82,83	0
5	SO4	D	709	5/5	0.91	0.34	-	68,76,78,87	0
6	EDO	C	709	4/4	0.93	0.20	-	30,34,35,35	0
2	NAG	B	704	14/15	0.86	0.26	-	59,67,70,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	SO4	B	707	5/5	0.96	0.28	-	73,74,81,82	0
5	SO4	A	707	5/5	0.87	0.26	-	65,69,76,83	0
2	NAG	A	704	14/15	0.80	0.36	-	59,71,84,84	0
2	NAG	C	701	14/15	0.70	0.40	-	75,92,97,100	0
2	NAG	D	701	14/15	0.81	0.38	-	78,83,85,86	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.