



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 2, 2023 – 04:38 PM EDT

PDB ID : 3WF0  
Title : Crystal structure of human beta-galactosidase in complex with 6S-NBI-DGJ  
Authors : Suzuki, H.; Ohto, U.; Shimizu, T.  
Deposited on : 2013-07-16  
Resolution : 2.20 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 2.36  
buster-report : 1.1.7 (2018)  
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) : 2.36

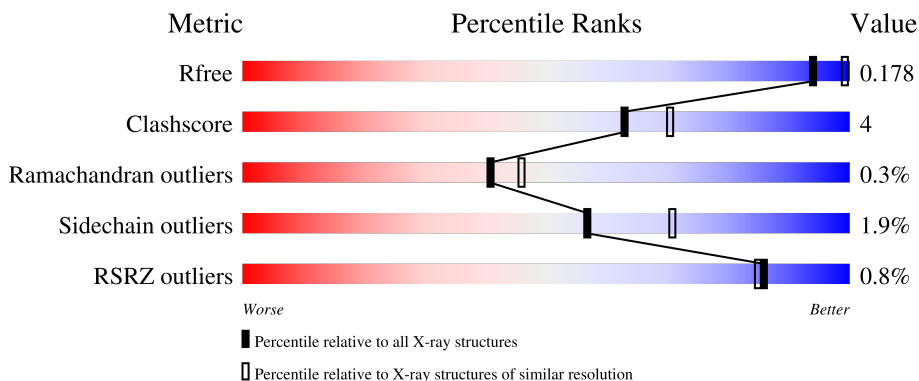
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	678	 80% 9% 11%
1	B	678	 78% 11% 11%
1	C	678	 79% 10% 11%
1	D	678	 78% 11% 11%

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	D	703	-	-	-	X

## 2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 20914 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
			Total	C	N	O	S			
1	A	605	4842	3139	801	884	18	0	9	0
1	B	605	4829	3133	795	884	17	0	6	0
1	C	603	4820	3126	793	883	18	0	9	0
1	D	603	4811	3118	791	884	18	0	9	0

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 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

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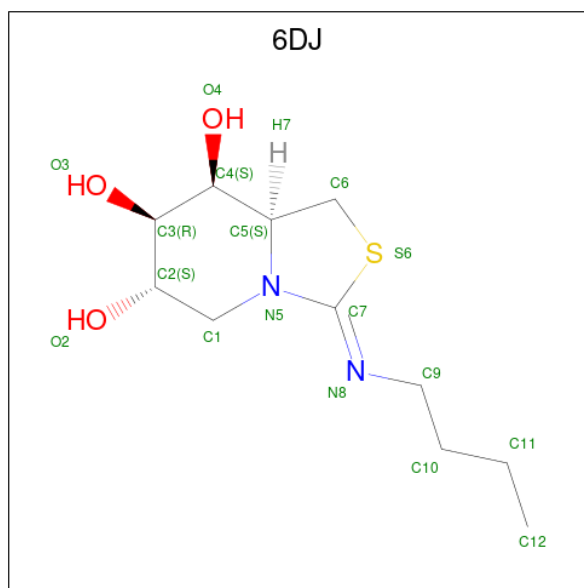
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		

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

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

- Molecule 4 is (3Z,6S,7R,8S,8aS)-3-(butylimino)hexahydro[1,3]thiazolo[3,4-a]pyridine-6,7,8-triol (three-letter code: 6DJ) (formula: C<sub>11</sub>H<sub>20</sub>N<sub>2</sub>O<sub>3</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			17	11	2	3	1		
4	B	1	Total	C	N	O	S	0	0
			17	11	2	3	1		
4	C	1	Total	C	N	O	S	0	0
			17	11	2	3	1		

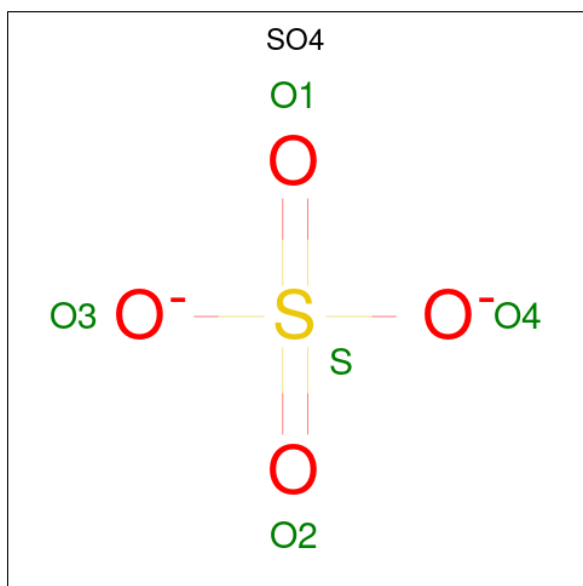
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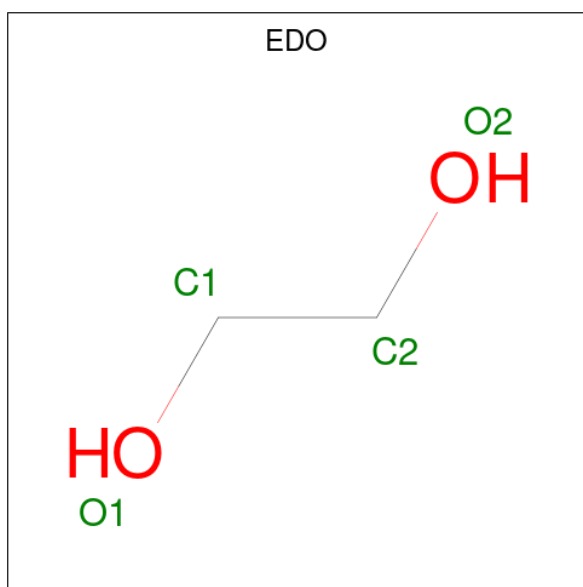
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	D	1	17	11	2	3	1	0	0

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



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

- 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 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0

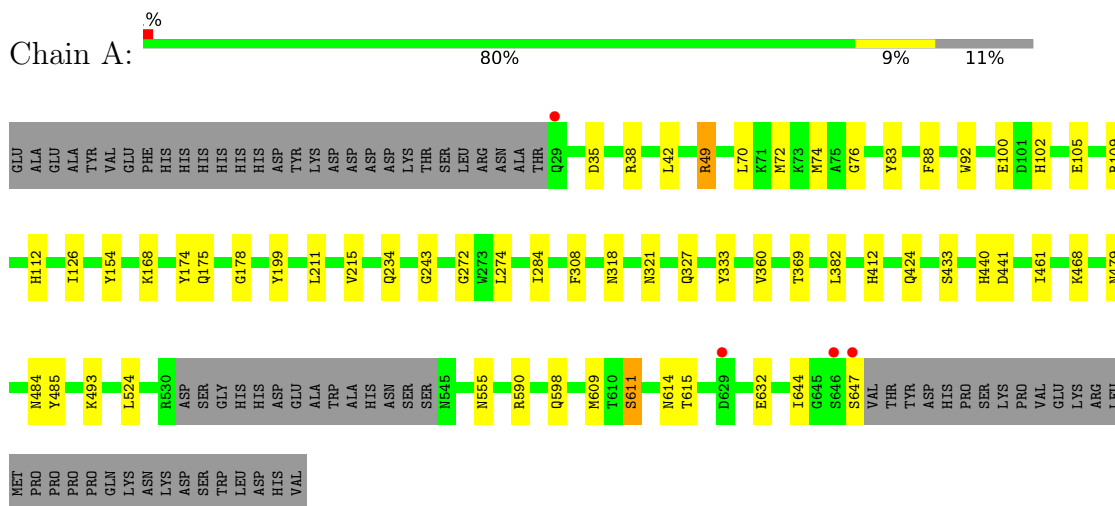
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	329	Total O 329 329	0	0
7	B	324	Total O 324 324	0	0
7	C	304	Total O 304 304	0	0
7	D	287	Total O 287 287	0	0

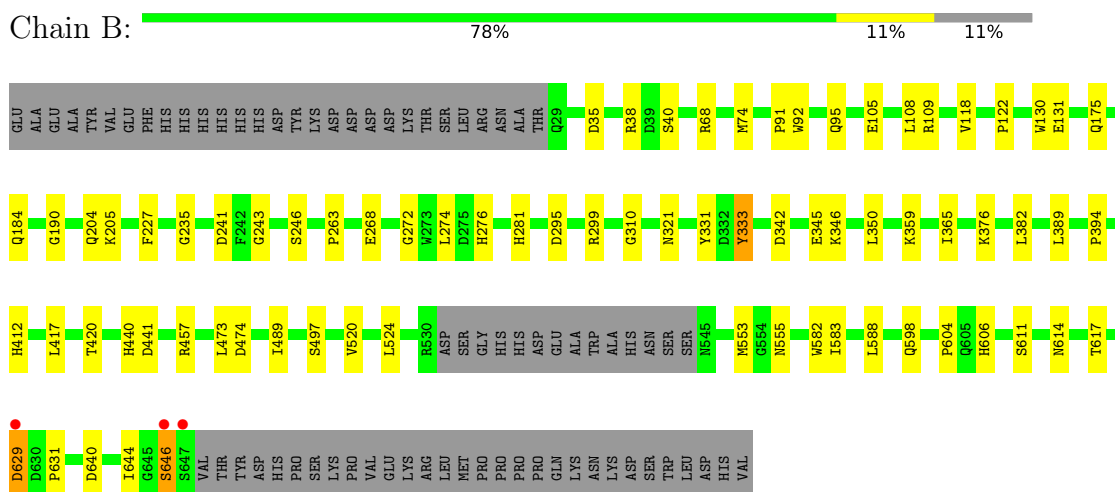
### 3 Residue-property plots [i](#)

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.

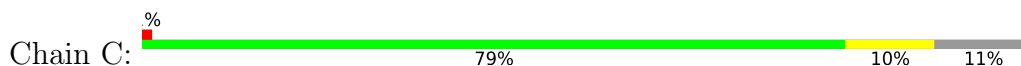
- Molecule 1: Beta-galactosidase

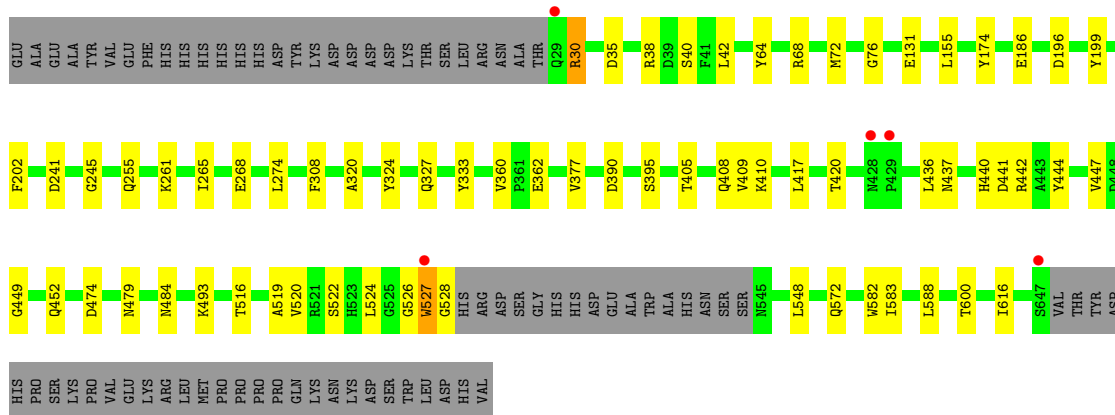


- Molecule 1: Beta-galactosidase

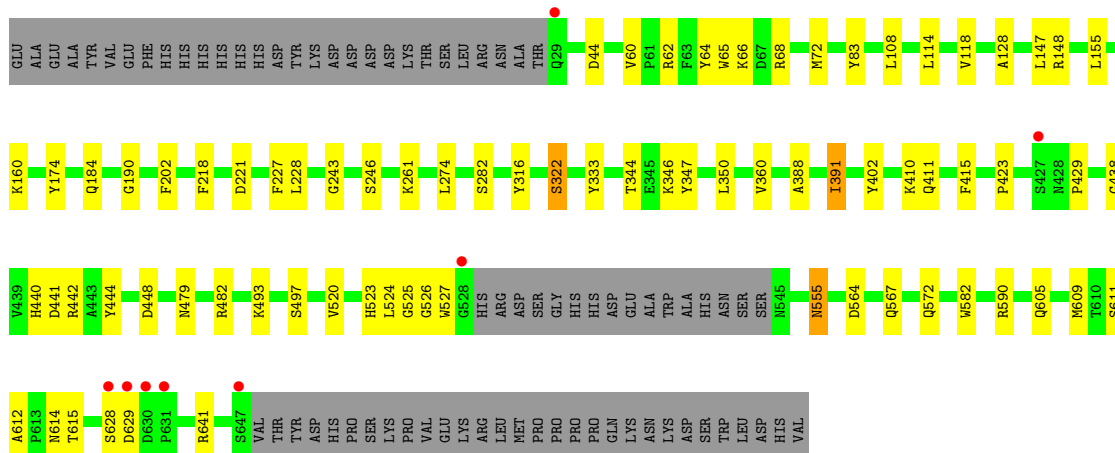
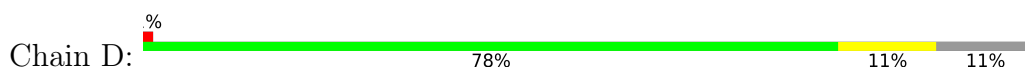


- Molecule 1: Beta-galactosidase





● Molecule 1: Beta-galactosidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.01Å 116.47Å 140.45Å 90.00° 92.23° 90.00°	Depositor
Resolution (Å)	43.41 – 2.20 43.41 – 2.20	Depositor EDS
% Data completeness (in resolution range)	90.1 (43.41-2.20) 90.1 (43.41-2.20)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.45 (at 2.20Å)	Xtrriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.177 , 0.230 0.181 , 0.178	Depositor DCC
$R_{free}$ test set	6997 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.0	Xtrriage
Anisotropy	0.038	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 33.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.023 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	20914	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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, 0.333 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: 6DJ, EDO, NAG, SO4, 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	0.67	0/5039	0.77	2/6872 (0.0%)
1	B	0.69	0/5010	0.76	4/6835 (0.1%)
1	C	0.67	0/5017	0.75	0/6844
1	D	0.67	0/5009	0.75	0/6834
All	All	0.68	0/20075	0.76	6/27385 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	68	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	A	49	ARG	CG-CD-NE	-5.39	100.48	111.80
1	B	342	ASP	CB-CG-OD1	5.36	123.12	118.30
1	A	168	LYS	CD-CE-NZ	-5.25	99.63	111.70
1	B	457	ARG	NE-CZ-NH2	-5.06	117.77	120.30
1	B	640	ASP	CB-CG-OD1	5.03	122.82	118.30

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	4842	0	4712	34	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	4829	0	4696	42	0
1	C	4820	0	4685	43	0
1	D	4811	0	4673	46	0
2	A	56	0	52	1	0
2	B	56	0	52	1	0
2	C	56	0	52	1	0
2	D	56	0	52	9	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	17	0	20	1	0
4	B	17	0	20	1	0
4	C	17	0	20	1	0
4	D	17	0	20	0	0
5	A	10	0	0	0	0
5	B	10	0	0	0	0
5	C	10	0	0	0	0
5	D	10	0	0	1	0
6	A	8	0	12	1	0
6	B	8	0	12	3	0
6	C	8	0	12	1	0
6	D	8	0	12	4	0
7	A	329	0	0	5	0
7	B	324	0	0	4	0
7	C	304	0	0	5	0
7	D	287	0	0	4	0
All	All	20914	0	19102	168	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 (168) 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:D:497:SER:HA	2:D:702:NAG:H81	1.10	1.10
1:D:497:SER:HA	2:D:702:NAG:C8	1.95	0.96
1:D:497:SER:CA	2:D:702:NAG:H81	2.03	0.86
1:C:405:THR:H	1:C:408:GLN:HE21	1.25	0.82
2:D:703:NAG:H83	2:D:703:NAG:H3	1.63	0.80
1:A:555:ASN:OD1	1:A:615:THR:HA	1.83	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:362:GLU:HG3	7:C:1046:HOH:O	1.85	0.75
1:A:327:GLN:HE22	1:A:484:ASN:HD21	1.37	0.71
1:A:318:ASN:HD21	1:A:590:ARG:HH21	1.36	0.71
1:D:402:TYR:CE1	2:D:702:NAG:H5	2.25	0.70
1:C:320:ALA:H	1:C:484:ASN:HD22	1.40	0.68
1:B:243:GLY:H	6:B:709:EDO:H22	1.58	0.67
1:C:35:ASP:HB2	1:C:42:LEU:HG	1.76	0.67
1:C:40:SER:HB2	7:C:813:HOH:O	1.95	0.67
1:D:243:GLY:H	6:D:709:EDO:H11	1.60	0.67
1:A:72:MET:HG2	1:A:308:PHE:CD2	2.30	0.67
1:B:382:LEU:HD23	1:B:524:LEU:HD12	1.79	0.65
1:C:245:GLY:H	6:C:709:EDO:H22	1.62	0.64
1:A:272:GLY:HA2	1:A:284:ILE:HG13	1.80	0.63
1:C:327:GLN:HE22	1:C:484:ASN:HD21	1.43	0.63
1:B:497:SER:HA	2:B:702:NAG:H81	1.80	0.62
1:A:112:HIS:HD2	7:A:846:HOH:O	1.83	0.60
1:D:388:ALA:O	1:D:391:ILE:HD12	2.01	0.60
1:C:390:ASP:HB2	7:C:1076:HOH:O	2.01	0.60
1:B:108:LEU:HD22	1:B:118:VAL:HG11	1.84	0.60
1:A:215:VAL:HG23	7:A:837:HOH:O	2.01	0.60
1:D:64:TYR:O	1:D:68:ARG:HG2	2.03	0.59
2:A:701:NAG:H81	7:A:1102:HOH:O	2.04	0.58
1:A:272:GLY:HA2	1:A:284:ILE:CG1	2.33	0.58
1:A:479:ASN:OD1	1:A:493:LYS:HE3	2.03	0.58
1:D:402:TYR:CZ	2:D:702:NAG:H5	2.39	0.57
1:A:88:PHE:O	1:A:102:HIS:HD2	1.86	0.57
1:D:60:VAL:HG11	1:D:68:ARG:HG3	1.85	0.56
1:B:611:SER:HB3	7:B:1093:HOH:O	2.05	0.56
1:C:30[A]:ARG:NH1	1:C:174:TYR:O	2.40	0.55
1:C:479:ASN:OD1	1:C:493:LYS:HE3	2.07	0.55
1:D:438:GLY:HA3	1:D:497:SER:OG	2.08	0.54
1:B:91:PRO:HD2	1:B:95:GLN:O	2.07	0.54
1:B:105:GLU:HB3	1:B:109:ARG:HH12	1.73	0.54
1:A:35:ASP:HB2	1:A:42:LEU:HG	1.90	0.53
1:C:255:GLN:HE22	1:C:265:ILE:H	1.57	0.53
1:B:241:ASP:HB2	1:B:268:GLU:HB2	1.90	0.53
1:D:322[A]:SER:OG	5:D:708:SO4:O4	2.21	0.52
1:A:112:HIS:CD2	7:A:846:HOH:O	2.61	0.52
1:A:105:GLU:O	1:A:109:ARG:HG3	2.10	0.51
1:D:246[B]:SER:OG	6:D:709:EDO:H12	2.11	0.51
1:D:360:VAL:HG22	7:D:1087:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:ARG:HG2	7:B:1096:HOH:O	2.10	0.51
1:C:527:TRP:O	1:C:528:GLY:C	2.49	0.51
1:A:382:LEU:HD23	1:A:524:LEU:HD12	1.92	0.51
1:B:346:LYS:O	1:B:350:LEU:HG	2.10	0.51
1:C:526:GLY:C	1:C:527:TRP:HE3	2.14	0.51
1:B:281:HIS:CE1	1:B:644:ILE:HD12	2.46	0.50
1:D:555:ASN:HB2	1:D:614:ASN:O	2.12	0.50
1:D:316:TYR:HD2	6:D:710:EDO:H22	1.75	0.50
1:A:174:TYR:HA	1:A:178:GLY:O	2.12	0.50
1:B:489:ILE:HD13	7:B:888:HOH:O	2.12	0.49
1:C:409:VAL:O	1:C:410:LYS:HB2	2.12	0.49
1:A:318:ASN:HD21	1:A:590:ARG:NH2	2.07	0.49
1:D:572:GLN:NE2	7:D:873:HOH:O	2.43	0.49
1:D:555:ASN:HD22	1:D:555:ASN:N	2.10	0.49
1:C:519:ALA:HB1	1:C:524:LEU:HD13	1.94	0.49
2:D:703:NAG:C8	2:D:703:NAG:C1	2.91	0.49
1:D:44:ASP:OD2	1:D:174:TYR:OH	2.28	0.48
1:D:66:LYS:HE3	1:D:114:LEU:HD21	1.95	0.48
1:A:243:GLY:H	6:A:709:EDO:H12	1.78	0.48
1:B:417:LEU:HD11	1:B:474:ASP:HB3	1.96	0.48
1:B:243:GLY:H	6:B:709:EDO:C2	2.24	0.47
1:D:523:HIS:HE1	1:D:615:THR:HG21	1.79	0.47
2:D:701:NAG:O7	2:D:701:NAG:O3	2.26	0.47
1:D:62:ARG:HA	1:D:65:TRP:CD2	2.50	0.47
1:A:92:TRP:CH2	1:A:412:HIS:CE1	3.03	0.47
1:D:590:ARG:HB3	7:D:844:HOH:O	2.14	0.47
1:A:611:SER:HB2	1:B:611:SER:O	2.15	0.46
1:D:83:TYR:CE2	1:D:128:ALA:HB2	2.50	0.46
1:D:411:GLN:NE2	1:D:415:PHE:O	2.45	0.46
1:C:572[B]:GLN:HB3	1:C:600:THR:HG22	1.96	0.46
1:C:268:GLU:OE1	4:C:706:6DJ:H5	2.15	0.46
1:A:211:LEU:HB3	1:A:215:VAL:HG21	1.97	0.46
1:A:321:ASN:ND2	1:A:327:GLN:HE21	2.13	0.46
1:C:320:ALA:N	1:C:484:ASN:HD22	2.11	0.46
1:B:268:GLU:OE1	4:B:706:6DJ:H5	2.15	0.46
1:A:440:HIS:HA	1:A:441:ASP:HA	1.60	0.46
1:B:629:ASP:O	1:B:631:PRO:HD3	2.15	0.46
1:A:609[A]:MET:HG2	1:A:614:ASN:OD1	2.16	0.46
1:C:377:VAL:HG11	1:C:616:ILE:HD12	1.98	0.45
1:B:190:GLY:HA3	1:B:227:PHE:O	2.16	0.45
1:D:520:VAL:HG11	1:D:582:TRP:CG	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:PRO:HD2	1:B:184:GLN:O	2.17	0.45
1:B:246:SER:OG	6:B:709:EDO:H21	2.17	0.45
1:C:35:ASP:OD2	1:C:38:ARG:HB2	2.17	0.45
1:C:442:ARG:HD3	1:C:444:TYR:OH	2.17	0.45
1:D:108:LEU:HD22	1:D:118:VAL:HG11	1.99	0.45
1:C:583:ILE:HD12	1:C:588:LEU:HD11	1.97	0.45
1:B:175:GLN:HB3	7:B:802:HOH:O	2.16	0.44
1:B:310:GLY:HA3	1:B:331:TYR:O	2.17	0.44
1:B:92:TRP:CH2	1:B:412:HIS:CE1	3.05	0.44
1:C:440:HIS:HA	1:C:441:ASP:HA	1.65	0.44
1:C:522:SER:HA	1:C:527:TRP:HB2	2.00	0.44
1:D:190:GLY:HA3	1:D:227:PHE:O	2.18	0.44
1:D:479:ASN:OD1	1:D:493:LYS:HE3	2.18	0.44
1:A:369:THR:HA	1:D:564[B]:ASP:OD2	2.18	0.44
1:B:40:SER:HA	1:B:263:PRO:HB3	1.99	0.44
1:D:440:HIS:HA	1:D:441:ASP:HA	1.77	0.44
1:B:553:MET:SD	1:B:617:THR:HG23	2.58	0.43
1:C:447:VAL:HG23	1:C:452:GLN:HG3	2.00	0.43
1:B:35:ASP:OD2	1:B:38:ARG:HB2	2.17	0.43
1:A:598:GLN:HA	1:A:644:ILE:HA	2.00	0.43
1:C:527:TRP:HE3	1:C:527:TRP:N	2.15	0.43
2:D:703:NAG:H83	2:D:703:NAG:C3	2.40	0.43
1:C:72:MET:HE3	1:C:308:PHE:HB3	2.01	0.43
1:A:92:TRP:CZ3	1:A:412:HIS:CE1	3.06	0.43
1:B:130:TRP:O	1:B:131:GLU:C	2.57	0.43
1:B:555:ASN:HA	1:B:614:ASN:O	2.19	0.43
1:B:420:THR:HG22	1:B:473:LEU:HD23	2.01	0.43
1:C:409:VAL:O	1:C:410:LYS:CB	2.67	0.43
1:D:62:ARG:HA	1:D:65:TRP:CE2	2.52	0.43
1:D:524:LEU:C	1:D:526:GLY:H	2.22	0.43
1:B:376:LYS:HE3	1:B:376:LYS:HB2	1.54	0.43
1:B:604:PRO:HB2	1:B:606:HIS:HD1	1.82	0.43
1:C:527:TRP:N	1:C:527:TRP:CE3	2.87	0.43
1:A:485:TYR:CE1	4:A:706:6DJ:H12	2.55	0.42
1:D:641:ARG:NH1	7:D:910:HOH:O	2.52	0.42
1:C:241:ASP:CB	1:C:268:GLU:HB2	2.49	0.42
1:C:520:VAL:HG11	1:C:582:TRP:CG	2.53	0.42
1:C:436:LEU:O	1:C:437:ASN:HB3	2.19	0.42
1:D:346:LYS:O	1:D:350:LEU:HG	2.20	0.42
1:B:295:ASP:O	1:B:299:ARG:HD2	2.19	0.42
1:C:241:ASP:HB2	1:C:268:GLU:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:GLN:HB3	7:A:805:HOH:O	2.18	0.42
1:A:433:SER:HA	1:A:461:ILE:O	2.19	0.42
1:C:68:ARG:O	1:C:72:MET:HG3	2.20	0.42
1:D:147:LEU:O	1:D:148:ARG:HB2	2.20	0.42
1:A:154:TYR:HE2	1:A:199:TYR:CE1	2.38	0.42
1:D:184:GLN:HB2	1:D:218:PHE:CZ	2.55	0.42
1:B:272:GLY:HA3	1:B:333:TYR:O	2.20	0.41
1:C:76:GLY:HA3	1:C:360:VAL:HG12	2.02	0.41
1:C:417:LEU:HD11	1:C:474:ASP:HB3	2.02	0.41
1:B:204:GLN:HE22	1:B:235:GLY:HA3	1.85	0.41
1:B:440:HIS:HA	1:B:441:ASP:HA	1.75	0.41
1:D:442:ARG:HD3	1:D:444:TYR:OH	2.20	0.41
1:D:316:TYR:HD2	6:D:710:EDO:C2	2.33	0.41
1:B:583:ILE:HD12	1:B:588:LEU:HD11	2.02	0.41
1:C:196:ASP:O	1:C:199:TYR:HB3	2.20	0.41
1:C:35:ASP:N	1:C:40:SER:O	2.47	0.41
1:B:276:HIS:NE2	1:B:321:ASN:ND2	2.60	0.41
1:D:221:ASP:HB2	1:D:228:LEU:HD23	2.03	0.41
1:A:83:TYR:CD1	1:A:126:ILE:HB	2.56	0.41
1:C:155:LEU:HD22	1:C:202:PHE:CD2	2.56	0.41
1:C:449:GLY:O	1:C:548:LEU:HD11	2.20	0.41
2:C:701:NAG:O3	7:C:1009:HOH:O	2.21	0.41
1:B:74:MET:HB3	1:B:365:ILE:HG12	2.01	0.41
1:C:64:TYR:O	1:C:68:ARG:HG2	2.20	0.40
1:D:344:THR:O	1:D:347:TYR:HB3	2.21	0.40
1:A:70:LEU:O	1:A:74:MET:HG2	2.21	0.40
1:A:76:GLY:HA3	1:A:360:VAL:HG22	2.02	0.40
1:B:598:GLN:HA	1:B:644:ILE:HA	2.03	0.40
1:B:520:VAL:HG11	1:B:582:TRP:CG	2.56	0.40
1:C:395:SER:HB2	7:C:914:HOH:O	2.21	0.40
1:D:482:ARG:HG2	1:D:493:LYS:HE2	2.04	0.40
1:D:609[A]:MET:SD	1:D:612:ALA:HB3	2.61	0.40
1:B:389:LEU:HD23	1:B:389:LEU:HA	1.88	0.40
1:B:394:PRO:HB3	1:D:609[A]:MET:HE1	2.03	0.40
1:D:68:ARG:O	1:D:72:MET:HG3	2.21	0.40
1:D:155:LEU:HD22	1:D:202:PHE:CD2	2.56	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	610/678 (90%)	585 (96%)	24 (4%)	1 (0%)	47	55
1	B	607/678 (90%)	581 (96%)	25 (4%)	1 (0%)	47	55
1	C	608/678 (90%)	580 (95%)	27 (4%)	1 (0%)	47	55
1	D	608/678 (90%)	583 (96%)	21 (4%)	4 (1%)	22	22
All	All	2433/2712 (90%)	2329 (96%)	97 (4%)	7 (0%)	41	46

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	646	SER
1	A	611	SER
1	D	628	SER
1	D	423	PRO
1	C	324	TYR
1	D	429	PRO
1	D	525	GLY

### 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	527/585 (90%)	517 (98%)	10 (2%)	57	71
1	B	524/585 (90%)	516 (98%)	8 (2%)	65	78
1	C	525/585 (90%)	515 (98%)	10 (2%)	57	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	525/585 (90%)	509 (97%)	16 (3%)	41	53
All	All	2101/2340 (90%)	2057 (98%)	44 (2%)	57	67

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38[A]	ARG
1	A	38[B]	ARG
1	A	49	ARG
1	A	100	GLU
1	A	234	GLN
1	A	274	LEU
1	A	333	TYR
1	A	424	GLN
1	A	468	LYS
1	A	647	SER
1	B	205	LYS
1	B	274	LEU
1	B	333	TYR
1	B	345[A]	GLU
1	B	345[B]	GLU
1	B	359	LYS
1	B	629	ASP
1	B	646	SER
1	C	30[A]	ARG
1	C	30[B]	ARG
1	C	131	GLU
1	C	186	GLU
1	C	261	LYS
1	C	274	LEU
1	C	333	TYR
1	C	420	THR
1	C	516	THR
1	C	527	TRP
1	D	160	LYS
1	D	261	LYS
1	D	274	LEU
1	D	282	SER
1	D	322[A]	SER
1	D	322[B]	SER
1	D	333	TYR
1	D	391	ILE

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Mol	Chain	Res	Type
1	D	410	LYS
1	D	448	ASP
1	D	527	TRP
1	D	555	ASN
1	D	605	GLN
1	D	611[A]	SER
1	D	611[B]	SER
1	D	629	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	102	HIS
1	A	112	HIS
1	A	318	ASN
1	A	321	ASN
1	A	355	GLN
1	A	484	ASN
1	B	102	HIS
1	B	204	GLN
1	B	321	ASN
1	B	355	GLN
1	B	412	HIS
1	C	255	GLN
1	C	408	GLN
1	C	440	HIS
1	C	484	ASN
1	D	97	GLN
1	D	424	GLN
1	D	523	HIS
1	D	572	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 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	D	702	1	14,14,15	0.29	0	17,19,21	0.62	0
2	NAG	C	704	1	14,14,15	0.30	0	17,19,21	0.61	0
4	6DJ	D	706	-	17,18,18	1.38	1 (5%)	17,25,25	2.16	6 (35%)
2	NAG	C	703	1	14,14,15	0.29	0	17,19,21	0.63	0
2	NAG	C	701	1	14,14,15	1.16	1 (7%)	17,19,21	1.74	4 (23%)
5	SO4	A	708	-	4,4,4	0.56	0	6,6,6	0.76	0
6	EDO	D	709	-	3,3,3	0.48	0	2,2,2	1.15	0
5	SO4	C	707	-	4,4,4	0.38	0	6,6,6	0.22	0
6	EDO	A	710	-	3,3,3	0.57	0	2,2,2	0.48	0
4	6DJ	A	706	-	17,18,18	1.03	1 (5%)	17,25,25	2.67	6 (35%)
6	EDO	B	710	-	3,3,3	0.54	0	2,2,2	0.40	0
2	NAG	B	703	1	14,14,15	0.86	0	17,19,21	1.99	4 (23%)
2	NAG	D	704	1	14,14,15	0.29	0	17,19,21	0.62	0
2	NAG	A	703	1	14,14,15	1.04	1 (7%)	17,19,21	1.96	5 (29%)
2	NAG	A	701	1	14,14,15	0.29	0	17,19,21	0.62	0
2	NAG	D	701	1	14,14,15	0.28	0	17,19,21	0.61	0
2	NAG	D	703	1	14,14,15	0.29	0	17,19,21	0.62	0
6	EDO	C	710	-	3,3,3	0.65	0	2,2,2	0.34	0
2	NAG	B	702	1	14,14,15	0.29	0	17,19,21	0.62	0
2	NAG	A	702	1	14,14,15	0.29	0	17,19,21	0.62	0
2	NAG	A	704	1	14,14,15	0.29	0	17,19,21	0.62	0
5	SO4	D	707	-	4,4,4	0.37	0	6,6,6	0.24	0
2	NAG	C	702	1	14,14,15	1.00	1 (7%)	17,19,21	0.94	0
5	SO4	B	708	-	4,4,4	0.54	0	6,6,6	1.20	1 (16%)
5	SO4	D	708	-	4,4,4	0.41	0	6,6,6	0.11	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	6DJ	B	706	-	17,18,18	1.13	1 (5%)	17,25,25	2.38	6 (35%)
5	SO4	C	708	-	4,4,4	0.47	0	6,6,6	0.75	0
6	EDO	A	709	-	3,3,3	0.61	0	2,2,2	0.26	0
5	SO4	B	707	-	4,4,4	0.40	0	6,6,6	0.26	0
2	NAG	B	704	1	14,14,15	0.29	0	17,19,21	0.62	0
4	6DJ	C	706	-	17,18,18	1.51	1 (5%)	17,25,25	2.51	5 (29%)
5	SO4	A	707	-	4,4,4	0.41	0	6,6,6	0.44	0
2	NAG	B	701	1	14,14,15	1.17	1 (7%)	17,19,21	1.91	5 (29%)
6	EDO	B	709	-	3,3,3	0.45	0	2,2,2	1.01	0
6	EDO	C	709	-	3,3,3	0.44	0	2,2,2	0.73	0
6	EDO	D	710	-	3,3,3	0.68	0	2,2,2	0.45	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	NAG	D	702	1	-	1/6/23/26	0/1/1/1
2	NAG	C	704	1	-	0/6/23/26	0/1/1/1
4	6DJ	D	706	-	-	1/5/34/34	0/2/2/2
2	NAG	C	703	1	-	1/6/23/26	0/1/1/1
2	NAG	C	701	1	-	0/6/23/26	0/1/1/1
6	EDO	D	709	-	-	1/1/1/1	-
6	EDO	A	710	-	-	1/1/1/1	-
4	6DJ	A	706	-	-	1/5/34/34	0/2/2/2
6	EDO	B	710	-	-	1/1/1/1	-
2	NAG	B	703	1	-	4/6/23/26	0/1/1/1
2	NAG	D	704	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	701	1	-	0/6/23/26	0/1/1/1
2	NAG	D	701	1	-	6/6/23/26	0/1/1/1
2	NAG	D	703	1	-	5/6/23/26	0/1/1/1
6	EDO	C	710	-	-	0/1/1/1	-
2	NAG	B	702	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	704	1	-	1/6/23/26	0/1/1/1
2	NAG	C	702	1	-	2/6/23/26	0/1/1/1
4	6DJ	B	706	-	-	1/5/34/34	0/2/2/2
6	EDO	A	709	-	-	0/1/1/1	-
2	NAG	B	704	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	6DJ	C	706	-	-	2/5/34/34	0/2/2/2
2	NAG	B	701	1	-	2/6/23/26	0/1/1/1
6	EDO	B	709	-	-	1/1/1/1	-
6	EDO	C	709	-	-	0/1/1/1	-
6	EDO	D	710	-	-	1/1/1/1	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	706	6DJ	C7-S6	-5.52	1.66	1.75
4	D	706	6DJ	C7-S6	-4.75	1.67	1.75
4	A	706	6DJ	C7-S6	-3.35	1.69	1.75
4	B	706	6DJ	C7-S6	-3.25	1.70	1.75
2	C	701	NAG	O5-C1	-2.47	1.39	1.43
2	B	701	NAG	O5-C1	-2.19	1.40	1.43
2	A	703	NAG	C2-N2	2.13	1.49	1.46
2	C	702	NAG	O5-C1	-2.03	1.40	1.43

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	706	6DJ	C9-N8-C7	7.03	129.30	117.84
2	B	703	NAG	C1-O5-C5	5.59	119.77	112.19
4	C	706	6DJ	C9-N8-C7	5.52	126.84	117.84
4	B	706	6DJ	C9-N8-C7	5.49	126.79	117.84
2	B	701	NAG	C1-O5-C5	5.28	119.35	112.19
4	C	706	6DJ	C10-C9-N8	-5.16	101.28	110.66
4	A	706	6DJ	C5-C6-S6	4.40	111.34	105.75
4	B	706	6DJ	C2-C1-N5	-4.31	103.59	110.30
4	A	706	6DJ	C1-N5-C5	-4.27	112.31	119.17
2	A	703	NAG	O7-C7-C8	-4.21	114.23	122.06
4	D	706	6DJ	C9-N8-C7	3.99	124.35	117.84
4	D	706	6DJ	C10-C9-N8	-3.88	103.61	110.66
2	A	703	NAG	C2-N2-C7	3.85	128.39	122.90
4	C	706	6DJ	C1-N5-C5	-3.82	113.04	119.17
4	D	706	6DJ	C2-C1-N5	-3.79	104.40	110.30
4	C	706	6DJ	C2-C1-N5	-3.75	104.47	110.30
4	B	706	6DJ	C5-C6-S6	3.68	110.43	105.75
4	A	706	6DJ	C10-C9-N8	-3.36	104.55	110.66
4	C	706	6DJ	C5-C6-S6	3.33	109.99	105.75
2	C	701	NAG	O4-C4-C3	-3.32	102.68	110.35
2	C	701	NAG	C4-C3-C2	3.26	115.79	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	706	6DJ	C1-N5-C5	-3.18	114.07	119.17
2	A	703	NAG	O7-C7-N2	3.12	127.69	121.95
2	B	701	NAG	C6-C5-C4	-3.03	105.90	113.00
2	C	701	NAG	C1-C2-N2	-3.02	105.34	110.49
2	A	703	NAG	O3-C3-C2	2.98	115.64	109.47
4	B	706	6DJ	O3-C3-C2	-2.98	104.29	109.99
4	A	706	6DJ	C2-C1-N5	-2.76	106.01	110.30
4	D	706	6DJ	C6-C5-C4	-2.74	111.58	115.72
2	B	703	NAG	O3-C3-C2	2.64	114.93	109.47
2	B	701	NAG	C8-C7-N2	-2.62	111.67	116.10
2	C	701	NAG	O3-C3-C4	-2.61	104.31	110.35
4	D	706	6DJ	C5-C6-S6	2.61	109.07	105.75
4	D	706	6DJ	C1-N5-C5	-2.56	115.05	119.17
2	B	701	NAG	O5-C5-C6	2.43	111.01	107.20
4	A	706	6DJ	C1-C2-C3	2.40	112.98	110.24
2	B	701	NAG	O7-C7-C8	2.25	126.23	122.06
2	A	703	NAG	O5-C1-C2	-2.18	107.84	111.29
2	B	703	NAG	O3-C3-C4	-2.16	105.36	110.35
4	B	706	6DJ	C10-C9-N8	-2.10	106.84	110.66
2	B	703	NAG	O5-C5-C4	2.10	115.94	110.83
5	B	708	SO4	O4-S-O3	-2.02	100.42	109.06

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	701	NAG	C8-C7-N2-C2
2	D	701	NAG	O7-C7-N2-C2
2	D	703	NAG	C3-C2-N2-C7
2	D	703	NAG	C8-C7-N2-C2
2	D	703	NAG	O7-C7-N2-C2
2	D	701	NAG	C1-C2-N2-C7
2	B	701	NAG	O5-C5-C6-O6
2	D	703	NAG	O5-C5-C6-O6
2	B	701	NAG	C4-C5-C6-O6
2	B	703	NAG	C4-C5-C6-O6
2	C	702	NAG	O5-C5-C6-O6
2	D	703	NAG	C4-C5-C6-O6
4	B	706	6DJ	C11-C10-C9-N8
4	C	706	6DJ	C11-C10-C9-N8
4	D	706	6DJ	C11-C10-C9-N8
2	B	703	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	C	702	NAG	C4-C5-C6-O6
2	D	701	NAG	O5-C5-C6-O6
2	D	701	NAG	C4-C5-C6-O6
6	A	710	EDO	O1-C1-C2-O2
2	B	703	NAG	C8-C7-N2-C2
6	B	709	EDO	O1-C1-C2-O2
6	D	709	EDO	O1-C1-C2-O2
6	D	710	EDO	O1-C1-C2-O2
4	A	706	6DJ	C9-C10-C11-C12
4	C	706	6DJ	C9-C10-C11-C12
2	B	703	NAG	O7-C7-N2-C2
2	D	701	NAG	C3-C2-N2-C7
2	A	704	NAG	C4-C5-C6-O6
2	C	703	NAG	O5-C5-C6-O6
6	B	710	EDO	O1-C1-C2-O2
2	D	702	NAG	C4-C5-C6-O6

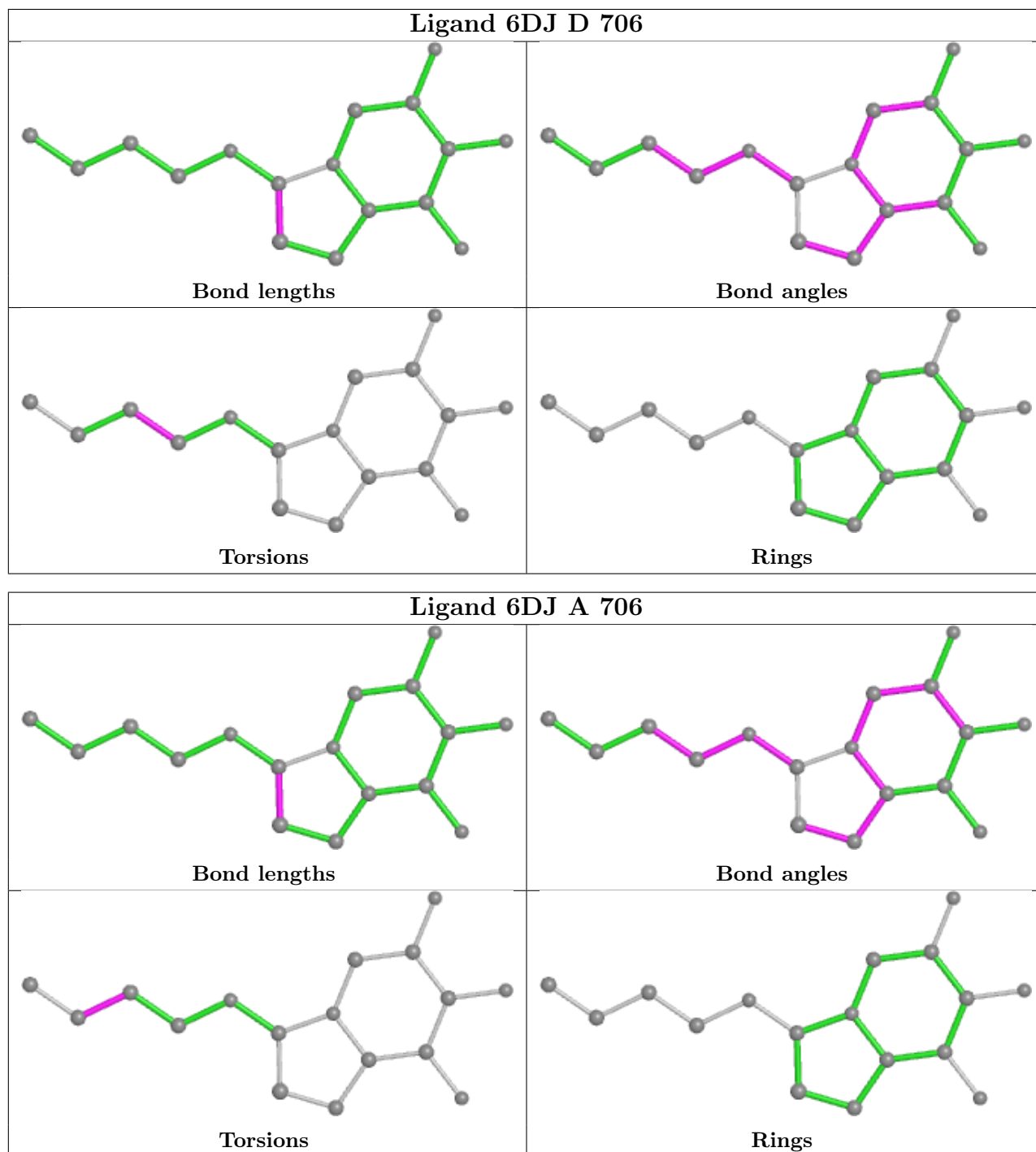
There are no ring outliers.

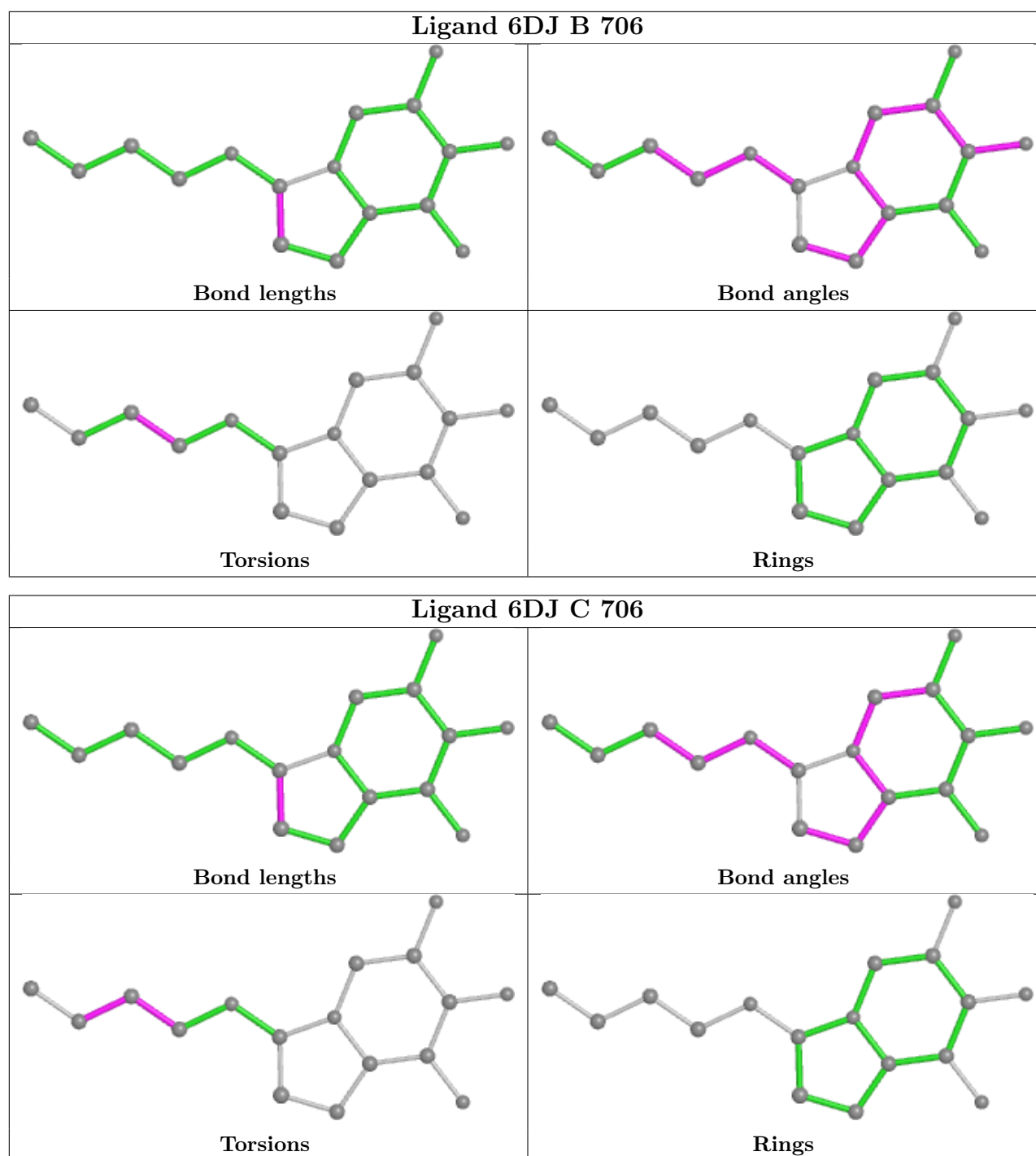
15 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	702	NAG	5	0
2	C	701	NAG	1	0
6	D	709	EDO	2	0
4	A	706	6DJ	1	0
2	A	701	NAG	1	0
2	D	701	NAG	1	0
2	D	703	NAG	3	0
2	B	702	NAG	1	0
5	D	708	SO4	1	0
4	B	706	6DJ	1	0
6	A	709	EDO	1	0
4	C	706	6DJ	1	0
6	B	709	EDO	3	0
6	C	709	EDO	1	0
6	D	710	EDO	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 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%)	-0.69	4 (0%) 87 86	10, 18, 34, 80	1 (0%)
1	B	605/678 (89%)	-0.71	3 (0%) 91 90	9, 17, 33, 76	0
1	C	603/678 (88%)	-0.63	5 (0%) 86 85	8, 19, 42, 74	0
1	D	603/678 (88%)	-0.57	8 (1%) 77 75	10, 20, 44, 82	0
All	All	2416/2712 (89%)	-0.65	20 (0%) 86 85	8, 18, 40, 82	1 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	646	SER	4.3
1	A	29	GLN	4.1
1	B	647	SER	3.7
1	D	628	SER	3.7
1	A	647	SER	3.4
1	A	646	SER	3.4
1	D	629	ASP	3.1
1	C	647	SER	3.1
1	A	629	ASP	2.6
1	D	29	GLN	2.6
1	B	629	ASP	2.5
1	D	528	GLY	2.5
1	C	29	GLN	2.4
1	C	429	PRO	2.2
1	D	427	SER	2.2
1	D	647	SER	2.1
1	D	631	PRO	2.1
1	D	630	ASP	2.1
1	C	527	TRP	2.0
1	C	428	ASN	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no 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.

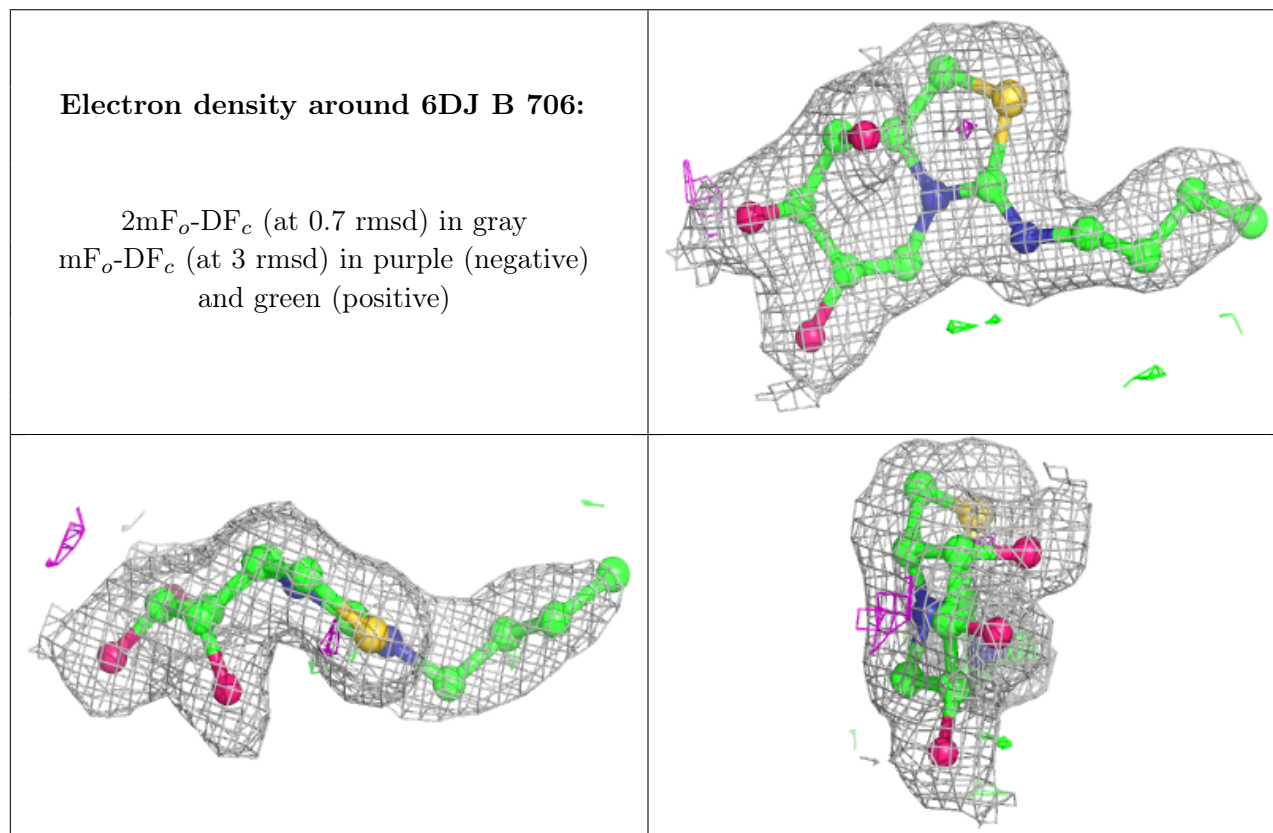
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	D	703	14/15	0.57	0.42	71,92,102,104	0
2	NAG	A	703	14/15	0.71	0.22	48,57,59,60	0
2	NAG	D	701	14/15	0.75	0.37	78,83,85,86	0
2	NAG	B	703	14/15	0.80	0.17	44,48,54,56	0
2	NAG	C	702	14/15	0.82	0.12	42,47,50,52	0
2	NAG	D	702	14/15	0.84	0.17	34,38,40,40	0
2	NAG	C	703	14/15	0.85	0.16	41,53,56,57	0
5	SO4	D	707	5/5	0.85	0.26	76,78,85,87	0
6	EDO	C	709	4/4	0.85	0.21	26,28,31,35	0
2	NAG	C	701	14/15	0.86	0.33	68,75,78,78	0
6	EDO	D	709	4/4	0.89	0.19	23,23,26,34	0
6	EDO	D	710	4/4	0.89	0.22	29,33,34,39	0
6	EDO	A	709	4/4	0.90	0.21	32,34,37,37	0
2	NAG	B	704	14/15	0.91	0.24	47,50,53,54	0
2	NAG	A	704	14/15	0.91	0.23	48,53,58,61	0
2	NAG	A	701	14/15	0.91	0.12	17,20,25,25	0
2	NAG	B	701	14/15	0.92	0.11	20,24,29,31	0
5	SO4	A	707	5/5	0.92	0.20	56,57,60,64	0
6	EDO	B	709	4/4	0.92	0.16	25,25,27,28	0
2	NAG	D	704	14/15	0.93	0.09	25,30,31,34	0
6	EDO	C	710	4/4	0.93	0.16	24,29,30,32	0
5	SO4	B	708	5/5	0.94	0.11	45,47,49,52	0
5	SO4	C	707	5/5	0.94	0.28	70,77,82,83	0
5	SO4	C	708	5/5	0.94	0.13	48,49,52,54	0
2	NAG	A	702	14/15	0.94	0.15	20,24,26,26	0
5	SO4	A	708	5/5	0.94	0.13	43,43,47,52	0
2	NAG	C	704	14/15	0.95	0.10	21,25,28,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	SO4	B	707	5/5	0.95	0.30	70,70,74,74	0
2	NAG	B	702	14/15	0.95	0.13	25,28,30,32	0
6	EDO	A	710	4/4	0.96	0.13	22,25,26,29	0
4	6DJ	B	706	17/17	0.96	0.13	14,16,31,33	0
4	6DJ	D	706	17/17	0.96	0.10	18,19,33,35	0
4	6DJ	A	706	17/17	0.97	0.11	18,19,30,31	0
4	6DJ	C	706	17/17	0.97	0.11	16,18,31,31	0
5	SO4	D	708	5/5	0.97	0.12	54,56,59,60	0
6	EDO	B	710	4/4	0.97	0.10	23,25,26,27	0
3	CL	B	705	1/1	0.99	0.10	16,16,16,16	0
3	CL	D	705	1/1	0.99	0.14	19,19,19,19	0
3	CL	C	705	1/1	1.00	0.09	14,14,14,14	0
3	CL	A	705	1/1	1.00	0.09	13,13,13,13	0

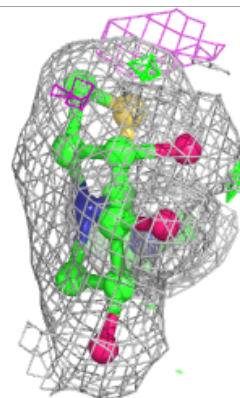
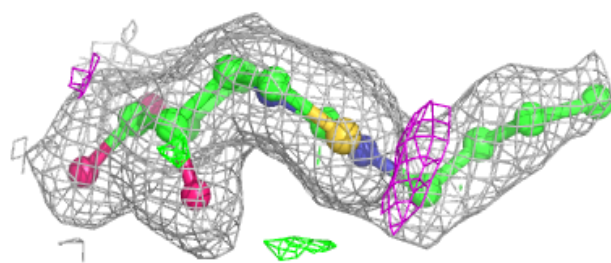
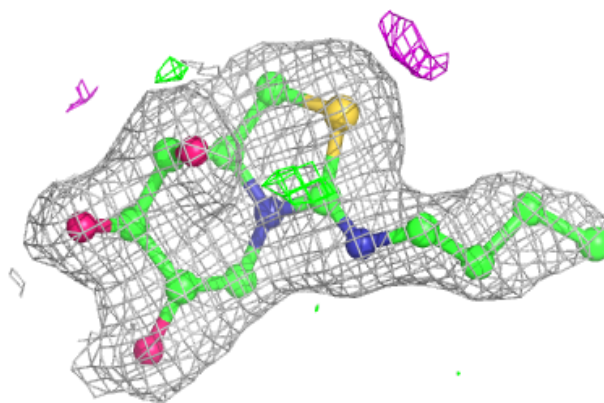
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



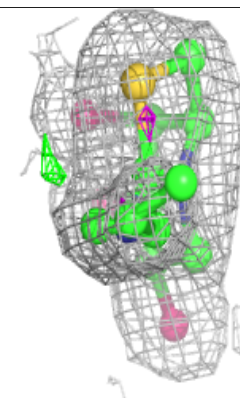
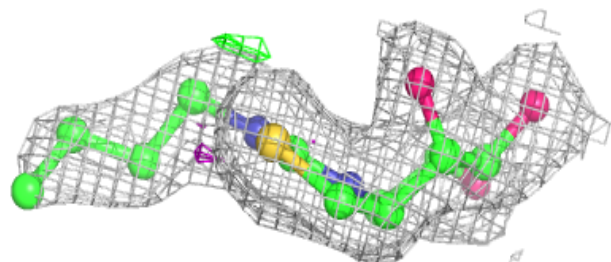
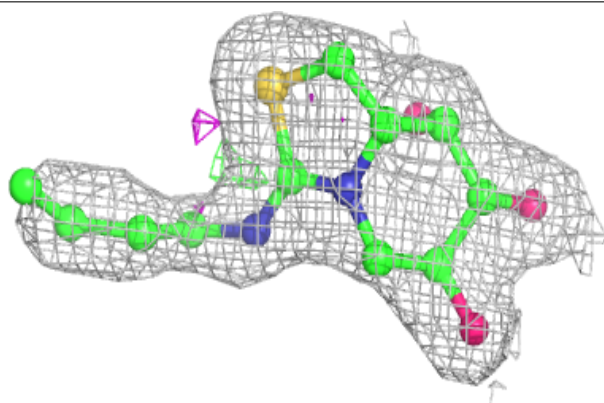


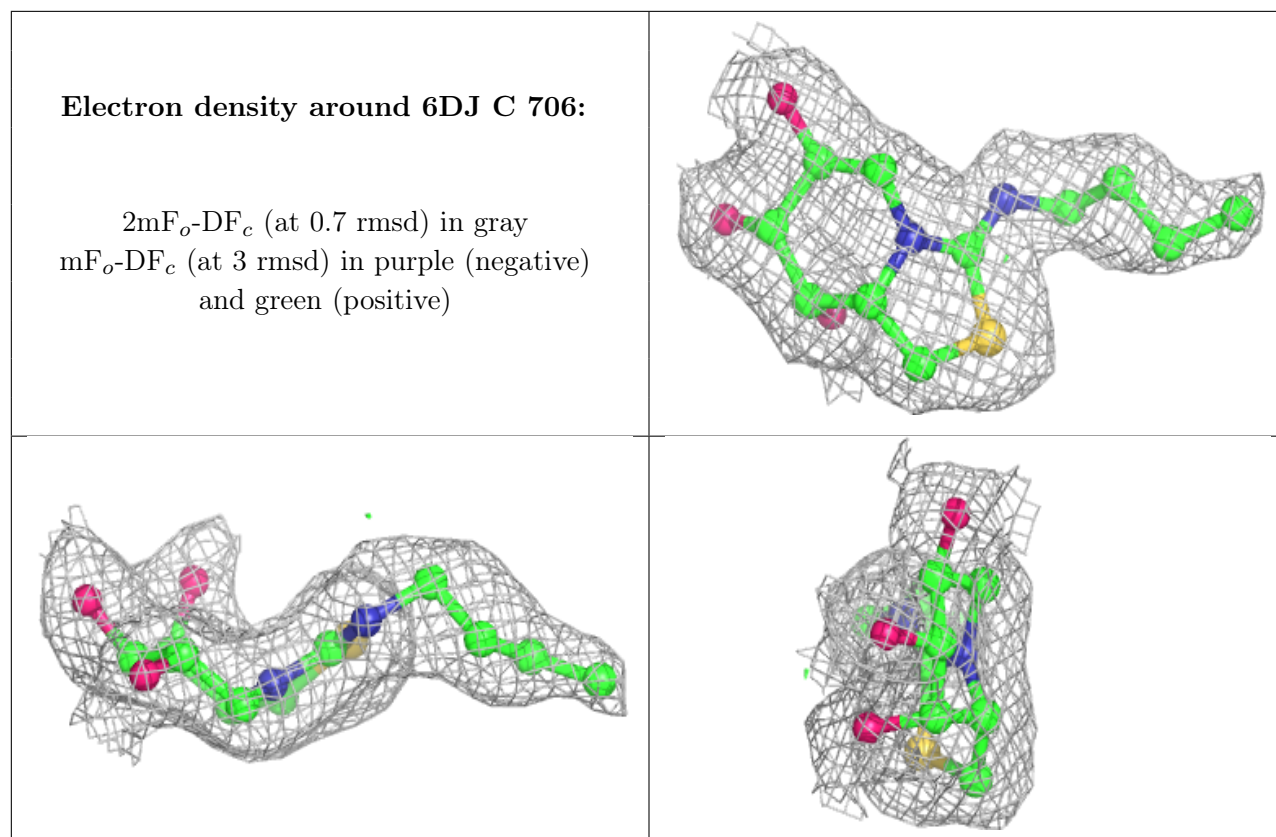
**Electron density around 6DJ D 706:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around 6DJ A 706:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 6.5 Other polymers [i](#)

There are no such residues in this entry.