



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:20 PM UTC

PDB ID : 3WF0 / pdb_00003wf0
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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

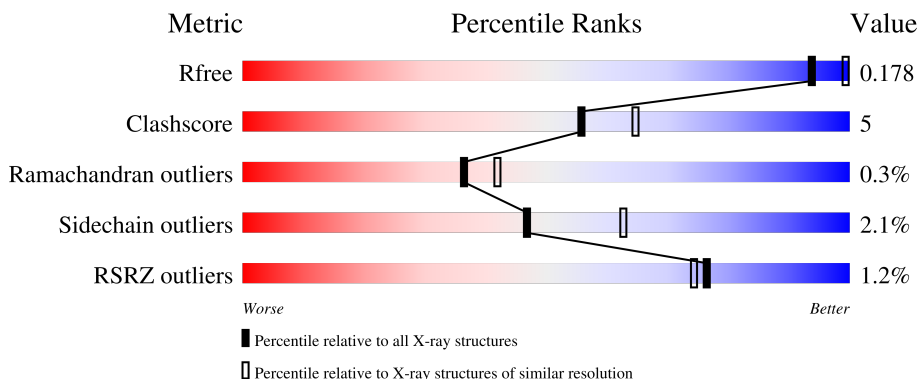
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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 ≥ 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	79% 10% 11%
1	B	678	77% 11% 11%
1	C	678	77% 11% 11%
1	D	678	77% 12% 11%

2 Entry composition

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
1	A	605	Total	C	N	O	S	0	9	0
			4842	3139	801	884	18			
1	B	605	Total	C	N	O	S	0	6	0
			4829	3133	795	884	17			
1	C	603	Total	C	N	O	S	0	9	0
			4820	3126	793	883	18			
1	D	603	Total	C	N	O	S	0	9	0
			4811	3118	791	884	18			

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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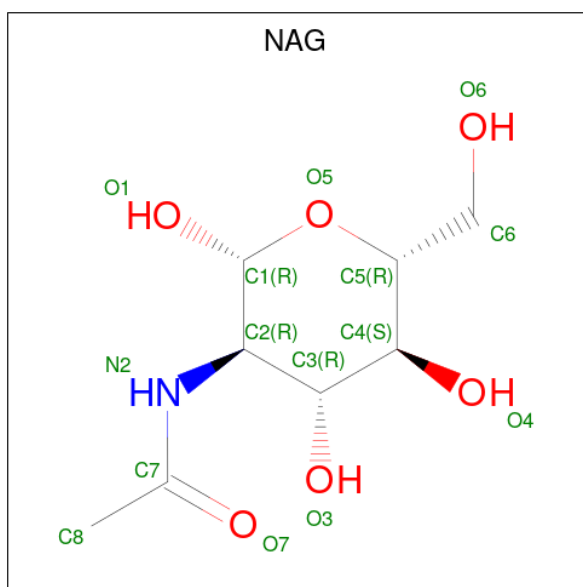
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	LYS	-	expression tag	UNP P16278
A	22	THR	-	expression tag	UNP P16278
A	23	SER	-	expression tag	UNP P16278
B	0	GLU	-	expression tag	UNP P16278
B	1	ALA	-	expression tag	UNP P16278
B	2	GLU	-	expression tag	UNP P16278
B	3	ALA	-	expression tag	UNP P16278
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 (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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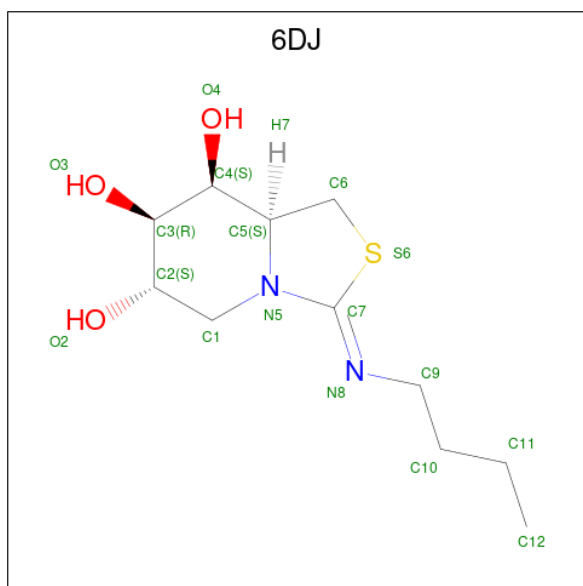
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is CHLORIDE ION (CCD ID: 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 (CCD ID: 6DJ) (formula: C₁₁H₂₀N₂O₃S).



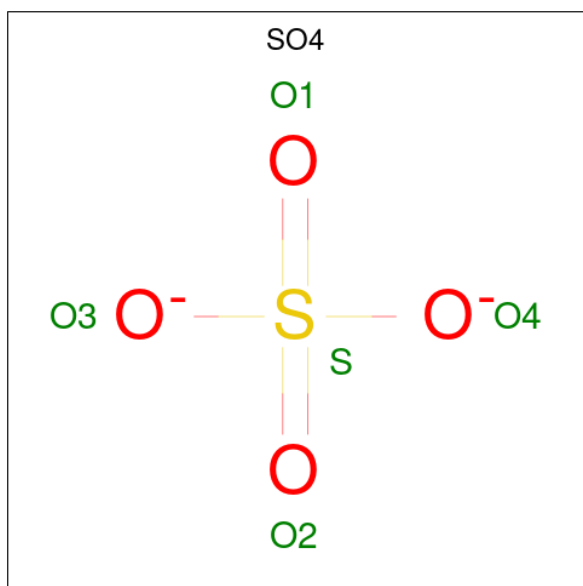
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			17	11	2	3	1		
4	B	1	Total	C	N	O	S	0	0
			17	11	2	3	1		
4	C	1	Total	C	N	O	S	0	0
			17	11	2	3	1		

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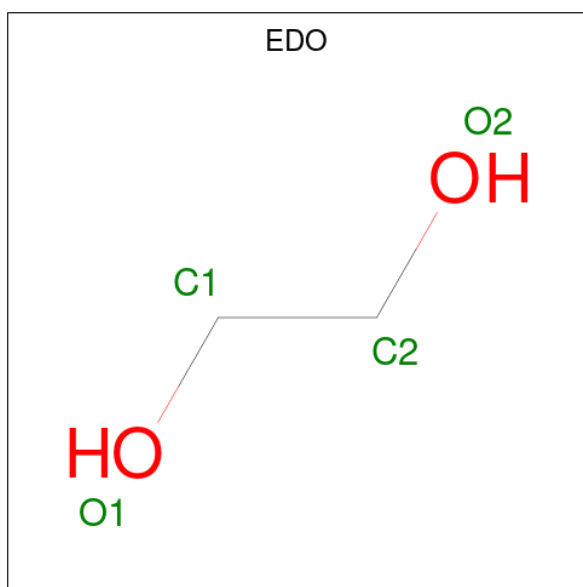
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	D	1	Total	C	N	O	S	0	0
			17	11	2	3	1		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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 (CCD ID: EDO) (formula: C₂H₆O₂).



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		

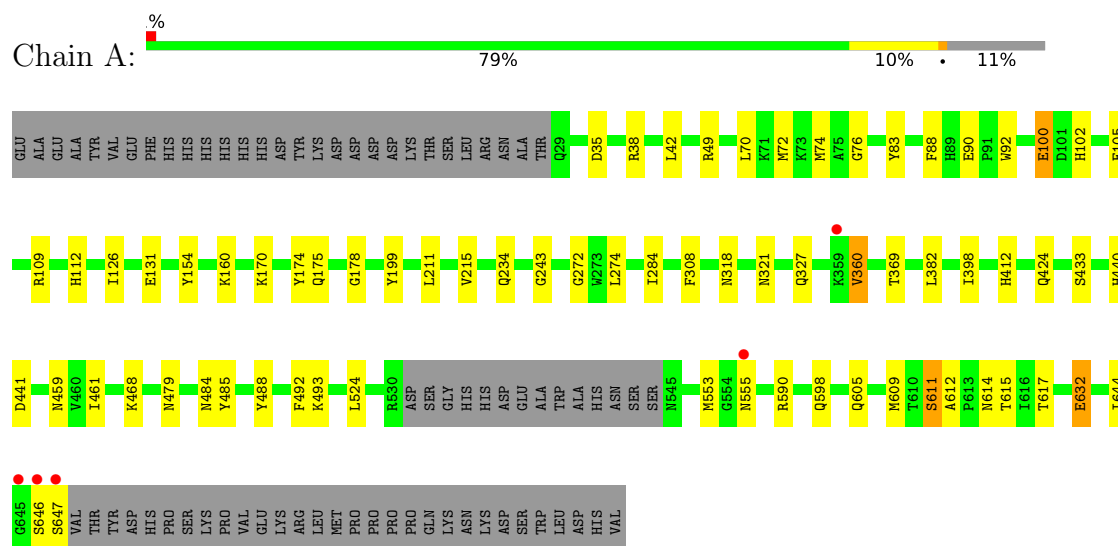
- Molecule 7 is water.

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

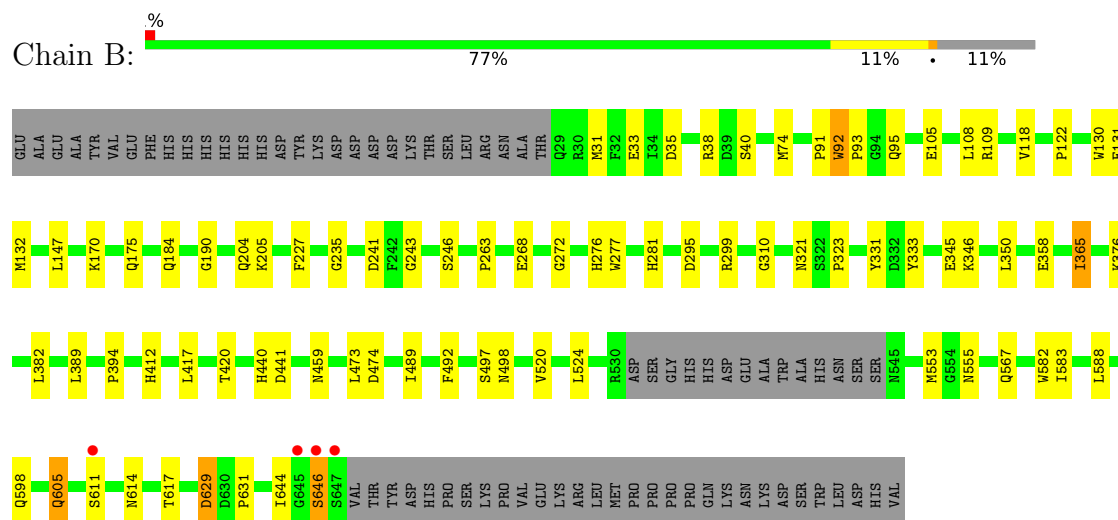
3 Residue-property plots

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.

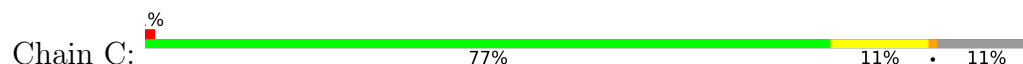
• Molecule 1: Beta-galactosidase

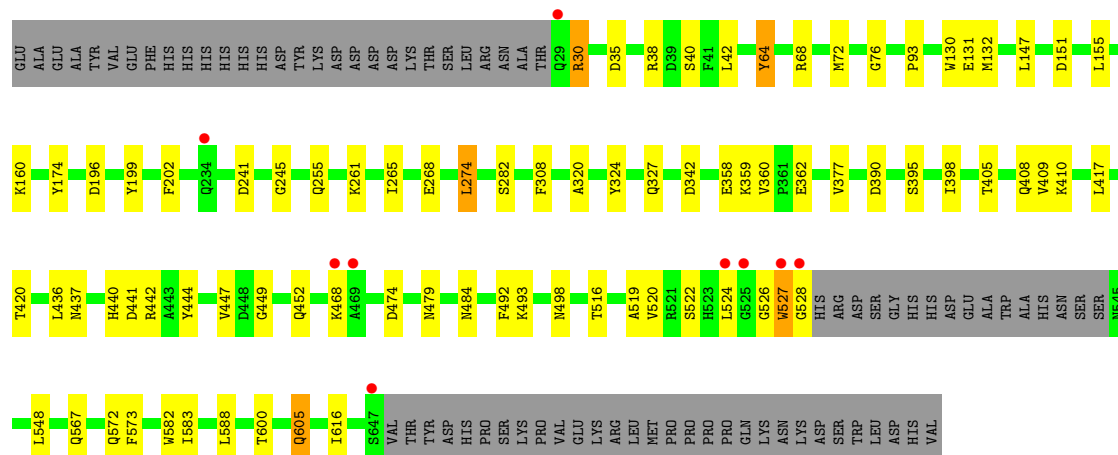


• Molecule 1: Beta-galactosidase

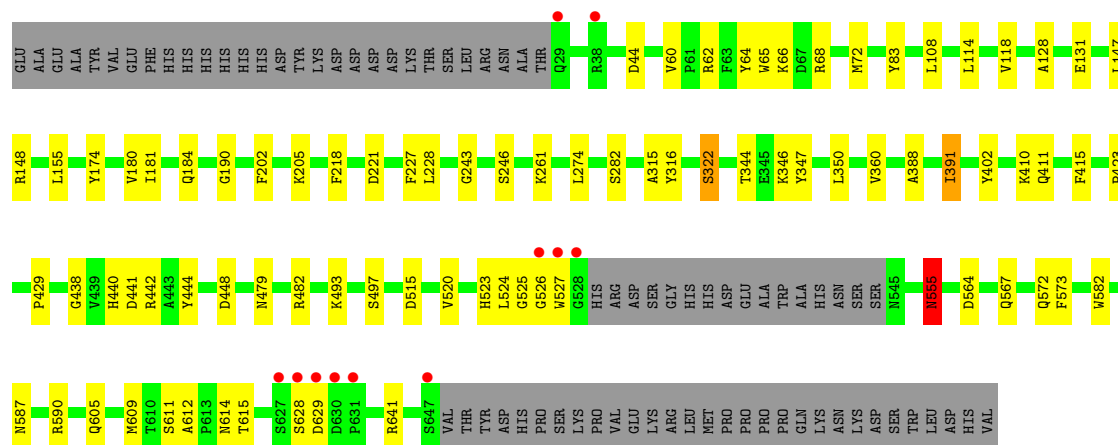
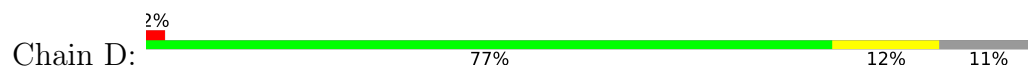


• Molecule 1: Beta-galactosidase





● Molecule 1: Beta-galactosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	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) 90.0 (43.41-2.20)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.20Å)	Xtriage
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 (4.52%)	wwPDB-VP
Wilson B-factor (Å ²)	20.0	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	20914	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage'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.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, NAG, EDO, CL, 6DJ

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	0.85	1/5039 (0.0%)	1.01	10/6872 (0.1%)
1	B	0.87	1/5010 (0.0%)	1.00	9/6835 (0.1%)
1	C	0.85	1/5017 (0.0%)	0.99	7/6844 (0.1%)
1	D	0.85	1/5009 (0.0%)	0.99	6/6834 (0.1%)
All	All	0.85	4/20075 (0.0%)	1.00	32/27385 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	492	PHE	C-O	5.39	1.30	1.23
1	A	492	PHE	C-O	5.30	1.30	1.24
1	D	555	ASN	C-O	-5.09	1.17	1.23
1	C	93	PRO	CA-C	5.06	1.58	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	646	SER	N-CA-C	9.31	116.41	108.78
1	B	459	ASN	N-CA-C	8.61	120.74	111.36
1	A	100	GLU	CB-CA-C	-7.50	107.93	116.63
1	D	555	ASN	N-CA-C	6.45	118.66	108.79
1	A	90	GLU	CA-C-N	6.32	126.52	119.32
1	A	90	GLU	C-N-CA	6.32	126.52	119.32
1	A	492	PHE	N-CA-C	6.22	117.72	111.07
1	D	573	PHE	CA-C-N	-6.20	113.37	119.76
1	D	573	PHE	C-N-CA	-6.20	113.37	119.76
1	C	573	PHE	CA-C-N	-6.18	113.31	119.99
1	C	573	PHE	C-N-CA	-6.18	113.31	119.99
1	A	488	TYR	N-CA-C	5.64	118.94	112.57
1	A	170	LYS	CA-C-N	-5.63	113.96	120.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	170	LYS	C-N-CA	-5.63	113.96	120.04
1	B	365	ILE	CA-C-N	-5.60	114.61	120.38
1	B	365	ILE	C-N-CA	-5.60	114.61	120.38
1	D	131	GLU	CB-CA-C	5.46	116.51	109.80
1	C	498	ASN	N-CA-C	5.40	118.11	110.50
1	B	170	LYS	CA-C-N	-5.36	114.30	119.87
1	B	170	LYS	C-N-CA	-5.36	114.30	119.87
1	C	64	TYR	N-CA-C	5.27	119.81	113.17
1	A	49	ARG	CG-CD-NE	-5.24	100.48	112.00
1	B	92	TRP	CA-C-N	-5.22	114.19	119.83
1	B	92	TRP	C-N-CA	-5.22	114.19	119.83
1	C	492	PHE	N-CA-C	-5.20	103.17	110.50
1	D	180	VAL	N-CA-C	5.19	115.08	108.12
1	D	181	ILE	N-CA-C	5.17	117.06	111.58
1	B	492	PHE	N-CA-C	-5.14	103.25	110.50
1	C	342	ASP	N-CA-C	5.13	117.33	110.35
1	B	498	ASN	N-CA-C	5.12	117.31	110.35
1	C	358	GLU	N-CA-C	5.09	116.27	108.42
1	A	459	ASN	N-CA-C	5.05	117.90	111.69

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	37	0
1	B	4829	0	4696	47	0
1	C	4820	0	4685	48	0
1	D	4811	0	4673	47	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	182	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 5.

All (182) 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
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:609[A]:MET:HE2	1:A:612:ALA:HB3	1.72	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:A:553:MET:HE2	1:A:617:THR:HG23	1.75	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:394:PRO:HB3	1:D:609[A]:MET:HE1	1.83	0.59
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:527:TRP:O	1:C:528:GLY:C	2.49	0.55
1:C:526:GLY:C	1:C:527:TRP:HE3	2.14	0.55
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:A:112:HIS:CD2	7:A:846:HOH:O	2.61	0.52
1:A:609[A]:MET:HE2	1:A:612:ALA:CB	2.40	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
1:B:299:ARG:HG2	7:B:1096:HOH:O	2.10	0.51
1:C:132:MET:HE3	1:C:147:LEU:HD11	1.92	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:322[A]:SER:OG	5:D:708:SO4:O4	2.21	0.48
1:A:243:GLY:H	6:A:709:EDO:H12	1.78	0.48
1:B:132:MET:HE3	1:B:147:LEU:HD11	1.95	0.48
1:B:417:LEU:HD11	1:B:474:ASP:HB3	1.96	0.48
1:D:524:LEU:C	1:D:526:GLY:H	2.22	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:B:130:TRP:O	1:B:131:GLU:C	2.57	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:C:72:MET:HE3	1:C:308:PHE:HB3	1.98	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:C:130:TRP:O	1:C:131:GLU:C	2.60	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:B:31:MET:HE2	1:B:33:GLU:OE2	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:A:92:TRP:CZ3	1:A:412:HIS:CE1	3.06	0.43
1:B:555:ASN:HA	1:B:614:ASN:O	2.19	0.43
1:D:315:ALA:HA	1:D:587:ASN:CG	2.43	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:B:376:LYS:HE3	1:B:376:LYS:HB2	1.54	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: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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:567:GLN:C	1:B:605:GLN:HB3	2.45	0.42
1:C:68:ARG:O	1:C:72:MET:HG3	2.20	0.42
1:C:274:LEU:O	1:C:282:SER:OG	2.36	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:C:151:ASP:C	1:C:151:ASP:OD1	2.62	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:C:436:LEU:O	1:C:437:ASN:HB3	2.19	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
1:C:567:GLN:C	1:C:605:GLN:HB3	2.45	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:B:276:HIS:O	1:B:277:TRP:C	2.63	0.40
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:92:TRP:O	1:B:93:PRO:C	2.63	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: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

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	610/678 (90%)	585 (96%)	24 (4%)	1 (0%)	43	51
1	B	607/678 (90%)	581 (96%)	25 (4%)	1 (0%)	43	51
1	C	608/678 (90%)	580 (95%)	27 (4%)	1 (0%)	43	51
1	D	608/678 (90%)	583 (96%)	21 (4%)	4 (1%)	18	19
All	All	2433/2712 (90%)	2329 (96%)	97 (4%)	7 (0%)	36	42

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

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%)	512 (97%)	15 (3%)	38	52
1	B	524/585 (90%)	517 (99%)	7 (1%)	61	76
1	C	525/585 (90%)	513 (98%)	12 (2%)	44	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	525/585 (90%)	509 (97%)	16 (3%)	36	49
All	All	2101/2340 (90%)	2051 (98%)	50 (2%)	47	58

All (50) 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	100	GLU
1	A	131	GLU
1	A	160	LYS
1	A	234	GLN
1	A	274	LEU
1	A	360	VAL
1	A	398	ILE
1	A	424	GLN
1	A	468	LYS
1	A	605	GLN
1	A	632[A]	GLU
1	A	632[B]	GLU
1	A	647	SER
1	B	205	LYS
1	B	345[A]	GLU
1	B	345[B]	GLU
1	B	358	GLU
1	B	605	GLN
1	B	629	ASP
1	B	646	SER
1	C	30[A]	ARG
1	C	30[B]	ARG
1	C	160	LYS
1	C	261	LYS
1	C	274	LEU
1	C	359	LYS
1	C	398	ILE
1	C	420	THR
1	C	468	LYS
1	C	516	THR
1	C	527	TRP
1	C	605	GLN
1	D	205	LYS
1	D	261	LYS

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Mol	Chain	Res	Type
1	D	274	LEU
1	D	282	SER
1	D	322[A]	SER
1	D	322[B]	SER
1	D	391	ILE
1	D	410	LYS
1	D	448	ASP
1	D	515	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 (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	81	GLN
1	A	102	HIS
1	A	112	HIS
1	A	175	GLN
1	A	318	ASN
1	A	321	ASN
1	A	355	GLN
1	A	484	ASN
1	A	605	GLN
1	B	102	HIS
1	B	204	GLN
1	B	321	ASN
1	B	352	ASN
1	B	412	HIS
1	C	112	HIS
1	C	204	GLN
1	C	255	GLN
1	C	408	GLN
1	C	484	ASN
1	D	204	GLN
1	D	355	GLN
1	D	523	HIS
1	D	572	GLN

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Mol	Chain	Res	Type
1	D	605	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 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	$\# Z > 2$
6	EDO	B	709	-	3,3,3	0.44	0	2,2,2	1.00	0
6	EDO	B	710	-	3,3,3	0.50	0	2,2,2	0.44	0
2	NAG	C	704	1	14,14,15	0.30	0	17,19,21	0.56	0
2	NAG	D	704	1	14,14,15	0.28	0	17,19,21	0.57	0
2	NAG	B	703	1	14,14,15	0.85	0	17,19,21	2.01	4 (23%)
5	SO4	B	707	-	4,4,4	0.49	0	6,6,6	0.22	0
5	SO4	C	708	-	4,4,4	0.55	0	6,6,6	0.65	0
4	6DJ	A	706	-	17,18,18	1.07	1 (5%)	16,25,25	2.67	6 (37%)
6	EDO	D	709	-	3,3,3	0.47	0	2,2,2	1.13	0
2	NAG	D	703	1	14,14,15	0.29	0	17,19,21	0.56	0
2	NAG	B	702	1	14,14,15	0.29	0	17,19,21	0.57	0
2	NAG	C	702	1	14,14,15	0.99	0	17,19,21	0.91	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	A	704	1	14,14,15	0.29	0	17,19,21	0.56	0
6	EDO	A	709	-	3,3,3	0.58	0	2,2,2	0.29	0
5	SO4	D	707	-	4,4,4	0.46	0	6,6,6	0.25	0
2	NAG	A	703	1	14,14,15	1.06	1 (7%)	17,19,21	2.01	5 (29%)
5	SO4	A	708	-	4,4,4	0.64	0	6,6,6	0.66	0
2	NAG	B	704	1	14,14,15	0.28	0	17,19,21	0.56	0
4	6DJ	C	706	-	17,18,18	1.58	1 (5%)	16,25,25	2.55	5 (31%)
2	NAG	D	702	1	14,14,15	0.29	0	17,19,21	0.56	0
2	NAG	A	701	1	14,14,15	0.29	0	17,19,21	0.56	0
5	SO4	A	707	-	4,4,4	0.50	0	6,6,6	0.43	0
2	NAG	B	701	1	14,14,15	1.15	1 (7%)	17,19,21	1.87	4 (23%)
4	6DJ	D	706	-	17,18,18	1.42	1 (5%)	16,25,25	2.09	4 (25%)
6	EDO	A	710	-	3,3,3	0.52	0	2,2,2	0.49	0
6	EDO	C	709	-	3,3,3	0.42	0	2,2,2	0.75	0
5	SO4	C	707	-	4,4,4	0.48	0	6,6,6	0.23	0
6	EDO	C	710	-	3,3,3	0.60	0	2,2,2	0.36	0
2	NAG	C	703	1	14,14,15	0.29	0	17,19,21	0.57	0
2	NAG	A	702	1	14,14,15	0.29	0	17,19,21	0.56	0
2	NAG	C	701	1	14,14,15	1.16	1 (7%)	17,19,21	1.75	4 (23%)
4	6DJ	B	706	-	17,18,18	1.15	1 (5%)	16,25,25	2.34	6 (37%)
5	SO4	D	708	-	4,4,4	0.50	0	6,6,6	0.12	0
6	EDO	D	710	-	3,3,3	0.65	0	2,2,2	0.42	0
5	SO4	B	708	-	4,4,4	0.62	0	6,6,6	1.04	0
2	NAG	D	701	1	14,14,15	0.28	0	17,19,21	0.56	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
6	EDO	B	709	-	-	1/1/1/1	-
6	EDO	B	710	-	-	1/1/1/1	-
2	NAG	C	704	1	-	0/6/23/26	0/1/1/1
2	NAG	D	704	1	-	0/6/23/26	0/1/1/1
2	NAG	B	703	1	-	4/6/23/26	0/1/1/1
4	6DJ	A	706	-	-	1/5/34/34	0/2/2/2
6	EDO	D	709	-	-	1/1/1/1	-
2	NAG	D	703	1	-	5/6/23/26	0/1/1/1
2	NAG	B	702	1	-	1/6/23/26	0/1/1/1
2	NAG	C	702	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	704	1	-	1/6/23/26	0/1/1/1
6	EDO	A	709	-	-	0/1/1/1	-
2	NAG	A	703	1	-	0/6/23/26	0/1/1/1
2	NAG	B	704	1	-	0/6/23/26	0/1/1/1
4	6DJ	C	706	-	-	2/5/34/34	0/2/2/2
2	NAG	D	702	1	-	1/6/23/26	0/1/1/1
2	NAG	A	701	1	-	0/6/23/26	0/1/1/1
2	NAG	B	701	1	-	2/6/23/26	0/1/1/1
4	6DJ	D	706	-	-	2/5/34/34	0/2/2/2
6	EDO	A	710	-	-	1/1/1/1	-
6	EDO	C	709	-	-	0/1/1/1	-
6	EDO	C	710	-	-	0/1/1/1	-
2	NAG	C	703	1	-	1/6/23/26	0/1/1/1
2	NAG	A	702	1	-	0/6/23/26	0/1/1/1
2	NAG	C	701	1	-	0/6/23/26	0/1/1/1
4	6DJ	B	706	-	-	1/5/34/34	0/2/2/2
6	EDO	D	710	-	-	1/1/1/1	-
2	NAG	D	701	1	-	6/6/23/26	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	706	6DJ	C7-S6	-5.85	1.66	1.75
4	D	706	6DJ	C7-S6	-5.04	1.67	1.75
4	A	706	6DJ	C7-S6	-3.56	1.69	1.75
4	B	706	6DJ	C7-S6	-3.47	1.70	1.75
2	C	701	NAG	O5-C1	-2.34	1.39	1.43
2	A	703	NAG	C2-N2	2.22	1.49	1.46
2	B	701	NAG	O5-C1	-2.07	1.40	1.43

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	706	6DJ	C9-N8-C7	7.36	129.30	117.83
4	C	706	6DJ	C9-N8-C7	5.78	126.84	117.83
4	B	706	6DJ	C9-N8-C7	5.75	126.79	117.83
2	B	703	NAG	C1-O5-C5	5.66	119.77	112.19
2	B	701	NAG	C1-O5-C5	5.35	119.35	112.19
4	C	706	6DJ	C10-C9-N8	-5.30	101.28	110.67
2	A	703	NAG	O7-C7-C8	-4.39	114.23	122.05
4	B	706	6DJ	C2-C1-N5	-4.31	103.59	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	706	6DJ	C1-N5-C5	-4.28	112.31	119.08
4	D	706	6DJ	C9-N8-C7	4.18	124.35	117.83
2	A	703	NAG	C2-N2-C7	4.10	128.39	122.90
4	D	706	6DJ	C10-C9-N8	-3.99	103.61	110.67
4	C	706	6DJ	C1-N5-C5	-3.82	113.04	119.08
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	A	706	6DJ	C10-C9-N8	-3.46	104.55	110.67
2	C	701	NAG	O4-C4-C3	-3.27	102.68	110.38
2	C	701	NAG	C4-C3-C2	3.26	115.79	111.02
2	C	701	NAG	C1-C2-N2	-3.23	105.34	110.43
2	A	703	NAG	O7-C7-N2	3.23	127.69	121.98
4	B	706	6DJ	C1-N5-C5	-3.17	114.07	119.08
2	A	703	NAG	O3-C3-C2	3.00	115.64	109.40
2	B	701	NAG	C6-C5-C4	-2.90	105.90	113.02
4	B	706	6DJ	O3-C3-C2	-2.82	104.29	110.05
4	A	706	6DJ	C5-C6-S6	2.77	111.34	106.31
4	A	706	6DJ	C2-C1-N5	-2.76	106.01	110.30
2	B	701	NAG	C8-C7-N2	-2.69	111.67	116.12
2	B	703	NAG	O3-C3-C2	2.66	114.93	109.40
2	C	701	NAG	O3-C3-C4	-2.57	104.31	110.38
4	D	706	6DJ	C1-N5-C5	-2.55	115.05	119.08
4	A	706	6DJ	C1-C2-C3	2.37	112.98	110.17
2	B	701	NAG	O7-C7-C8	2.35	126.23	122.05
4	B	706	6DJ	C5-C6-S6	2.27	110.43	106.31
2	A	703	NAG	O5-C1-C2	-2.23	107.84	111.29
4	B	706	6DJ	C10-C9-N8	-2.17	106.84	110.67
2	B	703	NAG	O3-C3-C4	-2.13	105.36	110.38
2	B	703	NAG	O5-C5-C4	2.10	115.94	110.83
4	C	706	6DJ	C5-C6-S6	2.03	109.99	106.31

There are no chirality outliers.

All (34) 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	B	701	NAG	O5-C5-C6-O6
2	D	703	NAG	O5-C5-C6-O6

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

There are no ring outliers.

15 monomers are involved in 25 short contacts:

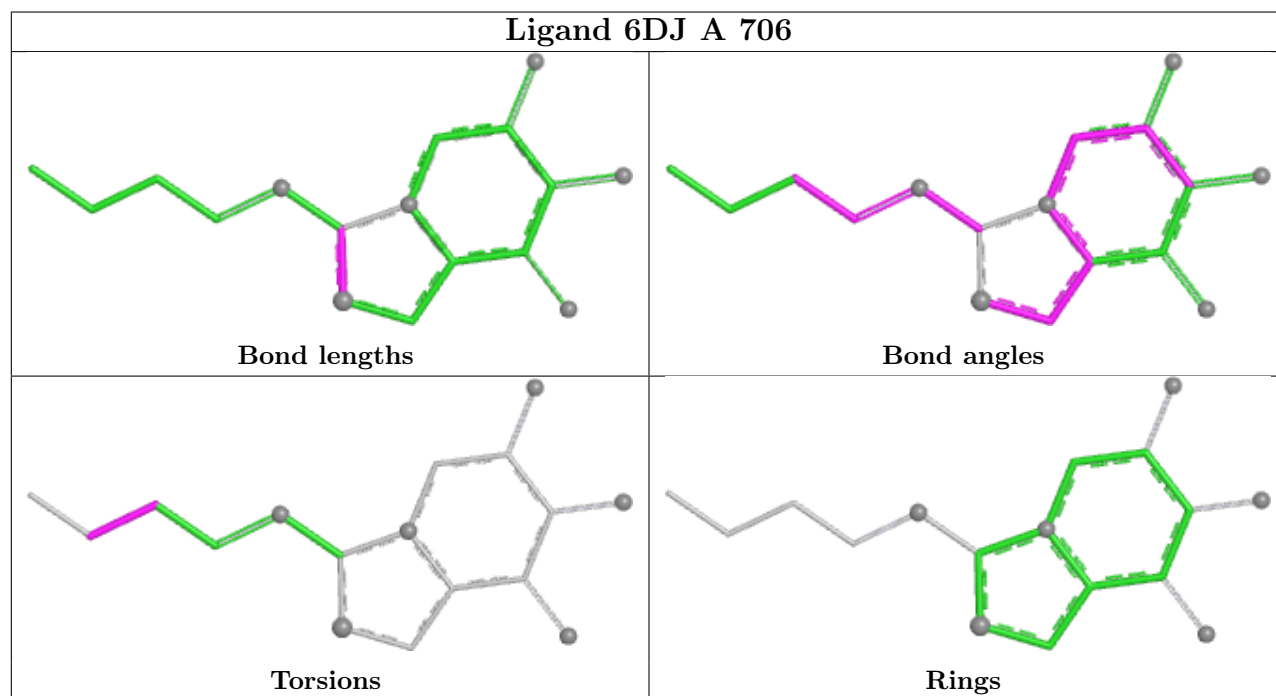
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	709	EDO	3	0
4	A	706	6DJ	1	0
6	D	709	EDO	2	0
2	D	703	NAG	3	0
2	B	702	NAG	1	0
6	A	709	EDO	1	0
4	C	706	6DJ	1	0
2	D	702	NAG	5	0
2	A	701	NAG	1	0
6	C	709	EDO	1	0

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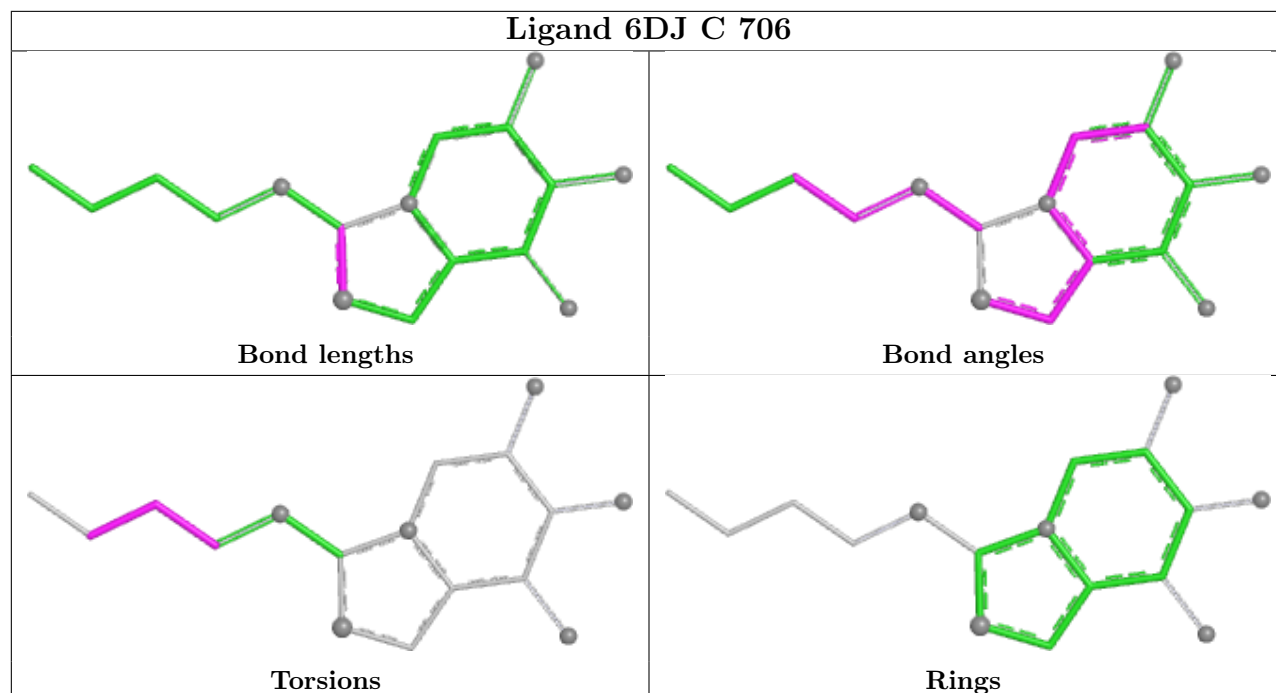
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	701	NAG	1	0
4	B	706	6DJ	1	0
5	D	708	SO4	1	0
6	D	710	EDO	2	0
2	D	701	NAG	1	0

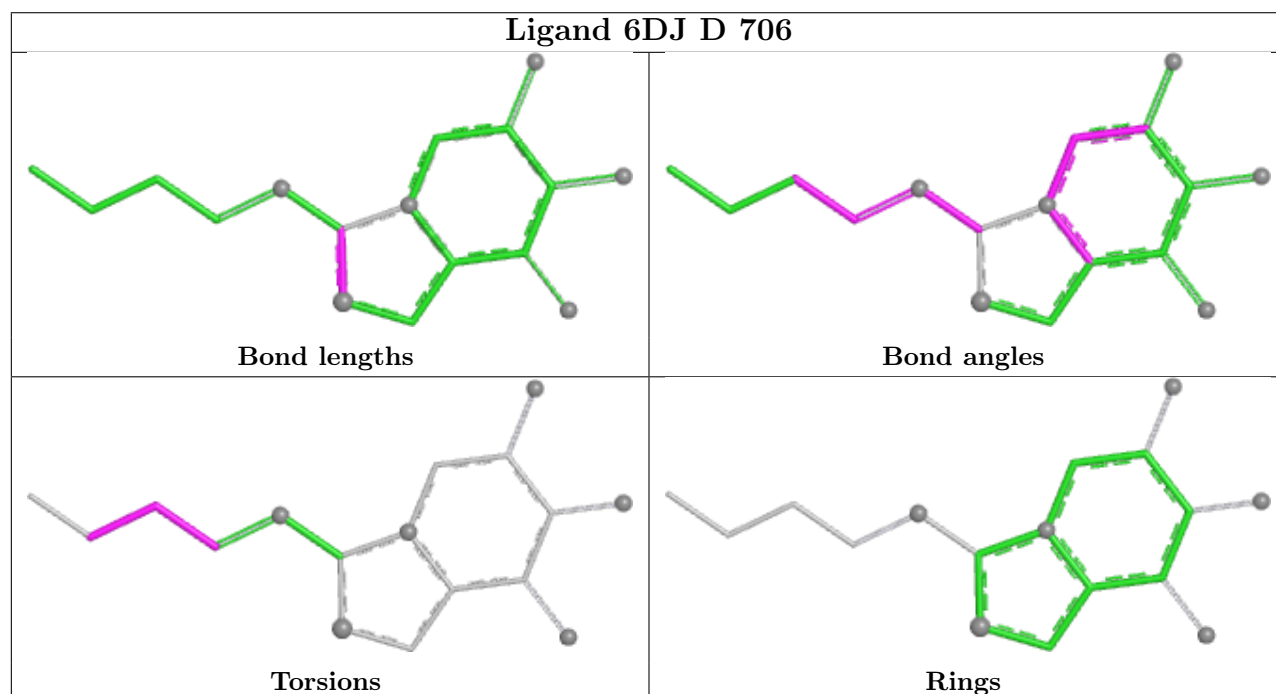
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

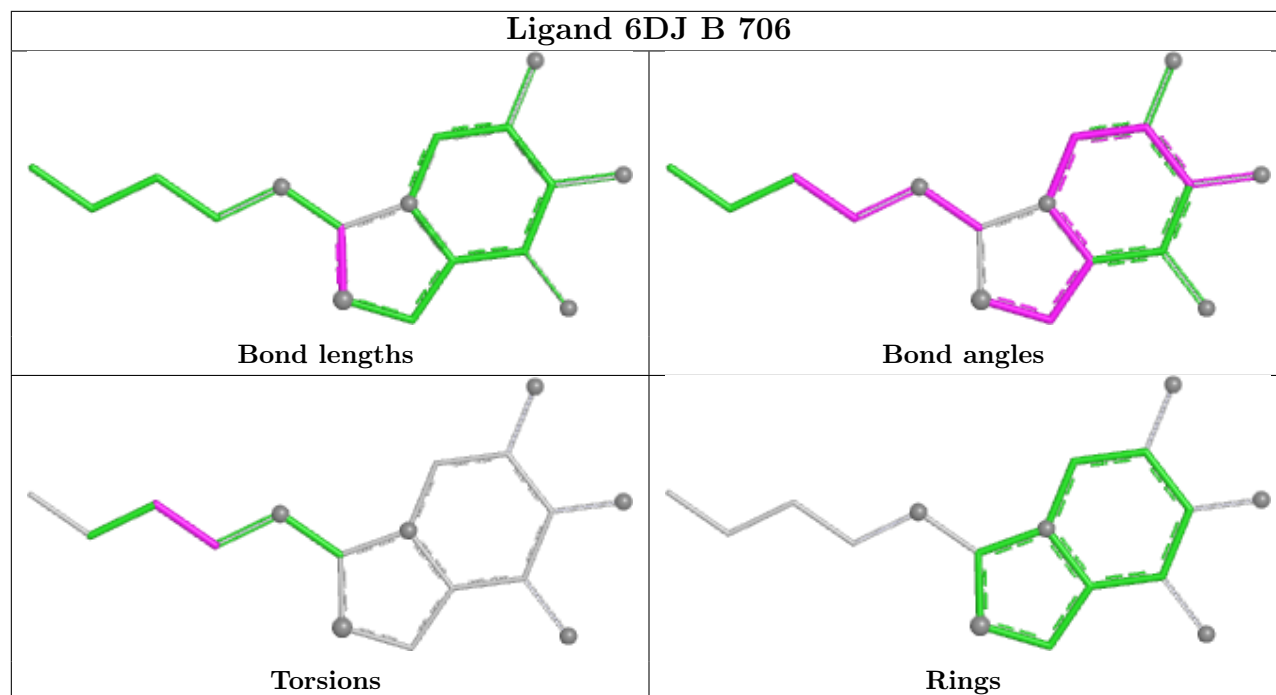


Ligand 6DJ C 706



Ligand 6DJ D 706





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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	605/678 (89%)	-0.45	5 (0%) 82 80	8, 18, 34, 80	9 (1%)
1	B	605/678 (89%)	-0.48	4 (0%) 84 82	9, 17, 33, 76	6 (0%)
1	C	603/678 (88%)	-0.35	9 (1%) 72 69	8, 19, 42, 74	9 (1%)
1	D	603/678 (88%)	-0.30	11 (1%) 67 64	10, 20, 44, 82	8 (1%)
All	All	2416/2712 (89%)	-0.39	29 (1%) 76 74	8, 18, 40, 82	32 (1%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	646	SER	5.3
1	C	527	TRP	5.2
1	A	645	GLY	4.5
1	B	647	SER	4.4
1	D	528	GLY	4.3
1	A	646	SER	4.1
1	A	647	SER	3.9
1	C	29	GLN	3.7
1	D	527	TRP	3.7
1	C	647	SER	3.3
1	D	647	SER	3.0
1	D	628	SER	2.9
1	A	555	ASN	2.9
1	B	611	SER	2.7
1	D	526	GLY	2.7
1	C	525	GLY	2.7
1	D	630	ASP	2.6
1	C	468	LYS	2.5
1	D	629	ASP	2.5
1	C	528	GLY	2.5
1	D	627	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	631	PRO	2.3
1	D	29	GLN	2.3
1	C	234	GLN	2.2
1	C	469	ALA	2.2
1	C	524	LEU	2.2
1	B	645	GLY	2.1
1	A	359	LYS	2.0
1	D	38	ARG	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 oligosaccharides 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	D	703	14/15	0.40	0.23	71,92,102,104	0
2	NAG	D	701	14/15	0.64	0.21	78,83,85,86	0
2	NAG	D	702	14/15	0.67	0.17	34,38,40,40	0
2	NAG	A	703	14/15	0.70	0.19	48,57,59,60	0
5	SO4	D	707	5/5	0.73	0.16	76,78,85,87	0
2	NAG	C	701	14/15	0.76	0.17	68,75,78,78	0
2	NAG	C	702	14/15	0.76	0.15	42,47,50,52	0
6	EDO	C	709	4/4	0.76	0.19	26,28,31,35	0
5	SO4	C	707	5/5	0.79	0.13	70,77,82,83	0
2	NAG	C	703	14/15	0.81	0.15	41,53,56,57	0
2	NAG	B	703	14/15	0.82	0.19	44,48,54,56	0
5	SO4	B	707	5/5	0.83	0.12	70,70,74,74	0
6	EDO	D	709	4/4	0.83	0.21	23,23,26,34	0
2	NAG	A	704	14/15	0.84	0.14	48,53,58,61	0
5	SO4	B	708	5/5	0.84	0.14	45,47,49,52	0

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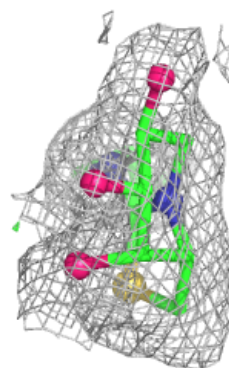
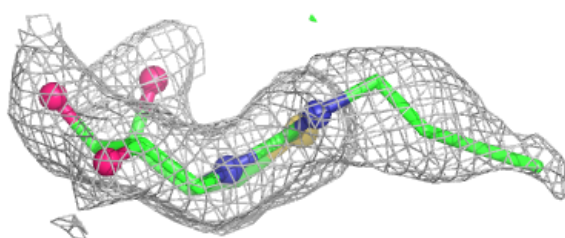
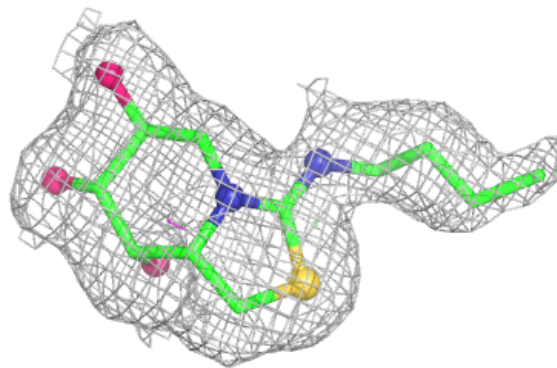
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	B	704	14/15	0.87	0.12	47,50,53,54	0
6	EDO	B	709	4/4	0.87	0.16	25,25,27,28	0
5	SO4	C	708	5/5	0.88	0.14	48,49,52,54	0
5	SO4	A	707	5/5	0.88	0.11	56,57,60,64	0
6	EDO	A	709	4/4	0.88	0.17	32,34,37,37	0
2	NAG	A	701	14/15	0.89	0.10	17,20,25,25	0
6	EDO	D	710	4/4	0.90	0.14	29,33,34,39	0
2	NAG	B	702	14/15	0.91	0.09	25,28,30,32	0
2	NAG	C	704	14/15	0.91	0.09	21,25,28,29	0
6	EDO	A	710	4/4	0.92	0.12	22,25,26,29	0
2	NAG	A	702	14/15	0.92	0.08	20,24,26,26	0
5	SO4	A	708	5/5	0.92	0.12	43,43,47,52	0
2	NAG	B	701	14/15	0.92	0.10	20,24,29,31	0
2	NAG	D	704	14/15	0.92	0.09	25,30,31,34	0
6	EDO	C	710	4/4	0.93	0.14	24,29,30,32	0
5	SO4	D	708	5/5	0.95	0.09	54,56,59,60	0
6	EDO	B	710	4/4	0.95	0.07	23,25,26,27	0
4	6DJ	C	706	17/17	0.96	0.07	16,18,31,31	0
4	6DJ	B	706	17/17	0.97	0.07	14,16,31,33	0
4	6DJ	D	706	17/17	0.97	0.07	18,19,33,35	0
4	6DJ	A	706	17/17	0.98	0.07	18,19,30,31	0
3	CL	B	705	1/1	0.99	0.03	16,16,16,16	0
3	CL	D	705	1/1	0.99	0.02	19,19,19,19	0
3	CL	C	705	1/1	1.00	0.04	14,14,14,14	0
3	CL	A	705	1/1	1.00	0.05	13,13,13,13	0

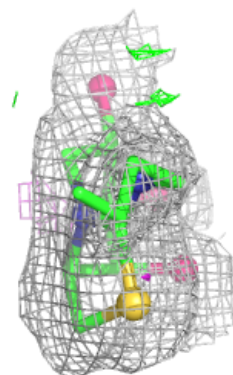
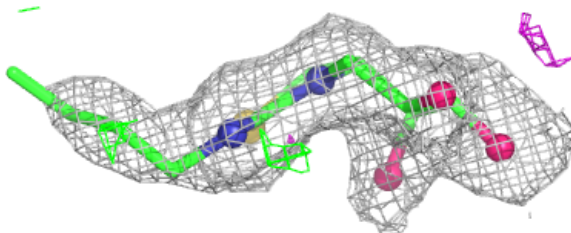
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 6DJ C 706:

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

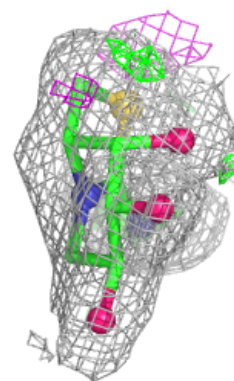
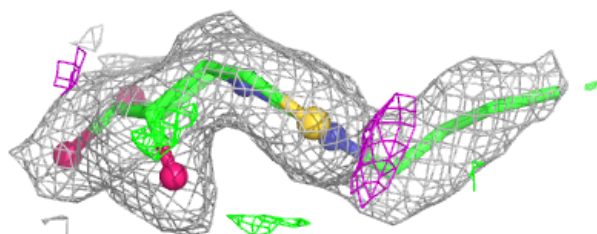
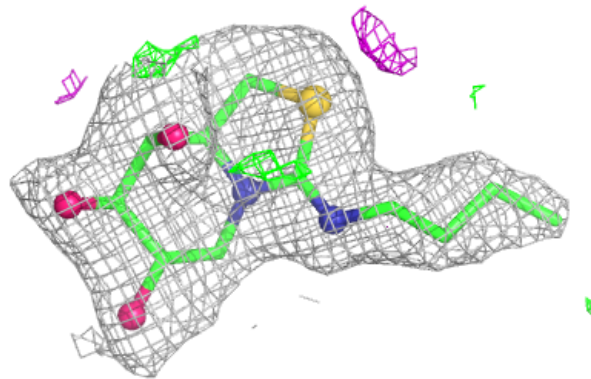
**Electron density around 6DJ B 706:**

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

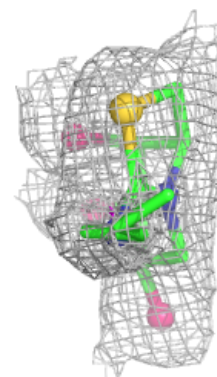
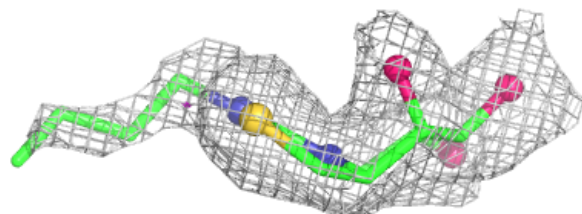
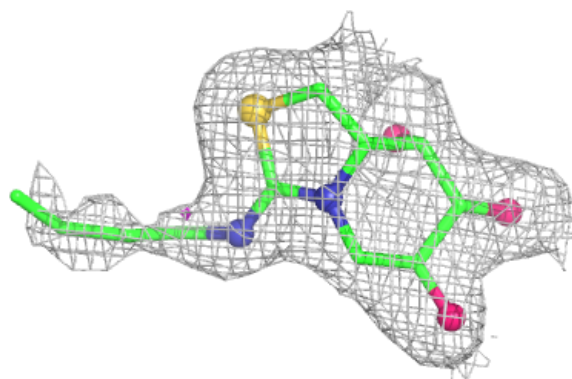


Electron density around 6DJ D 706:

$2mF_o-DF_c$ (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.