



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 23, 2022 – 12:08 PM EST

PDB ID : 7LVB  
Title : CHOLERA TOXIN B SUBUNIT WITH ATTACHED SIV EPITOPE  
Authors : Patskovsky, Y.; Cardozo, T.J.  
Deposited on : 2021-02-24  
Resolution : 2.25 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.26
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)	:	2.26

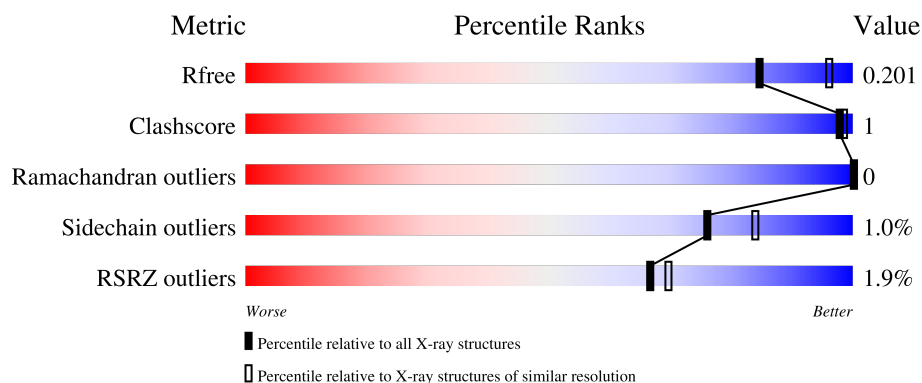
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	117	<div> <div>3%</div> <div>89%</div> <div>8%</div> </div>
1	B	117	<div> <div>2%</div> <div>91%</div> <div>8%</div> </div>
1	C	117	<div> <div>0%</div> <div>89%</div> <div>9%</div> </div>
1	D	117	<div> <div>2%</div> <div>90%</div> <div>9%</div> </div>
1	E	117	<div> <div>0%</div> <div>89%</div> <div>9%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	117	<div><div><div></div><div></div><div></div></div><div><div></div><div></div><div></div></div><div>3%</div><div>90%</div><div>• 8%</div></div>
1	G	117	<div><div><div></div><div></div><div></div></div><div><div></div><div></div><div></div></div><div>2%</div><div>90%</div><div>• 9%</div></div>
1	H	117	<div><div><div></div><div></div><div></div></div><div><div></div><div></div><div></div></div><div>2%</div><div>89%</div><div>• 9%</div></div>
1	I	117	<div><div><div></div><div></div><div></div></div><div><div></div><div></div><div></div></div><div>2%</div><div>88%</div><div>• • 9%</div></div>
1	J	117	<div><div><div></div><div></div><div></div></div><div><div></div><div></div><div></div></div><div>2%</div><div>90%</div><div>• • 9%</div></div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9406 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 Cholera enterotoxin B-subunit, Surface protein gp120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	107	Total	C	N	O	S	0	0	0
			847	533	148	161	5			
1	E	107	Total	C	N	O	S	0	0	0
			847	533	148	161	5			
1	F	108	Total	C	N	O	S	0	0	0
			856	538	149	164	5			
1	G	107	Total	C	N	O	S	0	0	0
			847	533	148	161	5			
1	H	107	Total	C	N	O	S	0	0	0
			847	533	148	161	5			
1	A	108	Total	C	N	O	S	0	0	0
			856	538	149	164	5			
1	B	108	Total	C	N	O	S	0	2	0
			868	546	151	166	5			
1	C	107	Total	C	N	O	S	0	2	0
			861	542	151	163	5			
1	I	107	Total	C	N	O	S	0	1	0
			854	538	150	161	5			
1	J	107	Total	C	N	O	S	0	0	0
			847	533	148	161	5			

There are 10 discrepancies between the modelled and reference sequences:

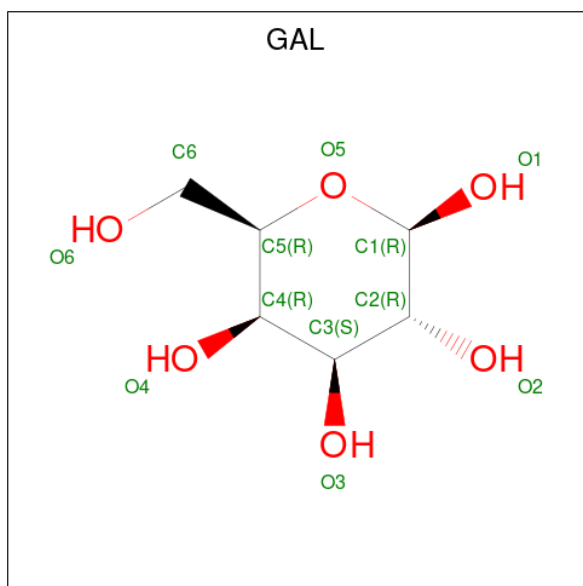
Chain	Residue	Modelled	Actual	Comment	Reference
D	107	ILE	ARG	variant	UNP P19503
E	107	ILE	ARG	variant	UNP P19503
F	107	ILE	ARG	variant	UNP P19503
G	107	ILE	ARG	variant	UNP P19503
H	107	ILE	ARG	variant	UNP P19503
A	107	ILE	ARG	variant	UNP P19503
B	107	ILE	ARG	variant	UNP P19503
C	107	ILE	ARG	variant	UNP P19503
I	107	ILE	ARG	variant	UNP P19503

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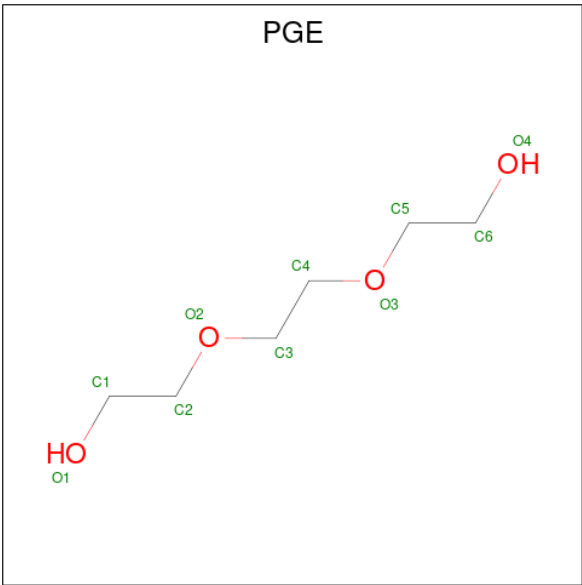
Chain	Residue	Modelled	Actual	Comment	Reference
J	107	ILE	ARG	variant	UNP P19503

- Molecule 2 is beta-D-galactopyranose (three-letter code: GAL) (formula:  $C_6H_{12}O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			12	6	6		
2	E	1	Total	C	O	0	0
			12	6	6		
2	F	1	Total	C	O	0	0
			12	6	6		
2	G	1	Total	C	O	0	0
			12	6	6		
2	H	1	Total	C	O	0	0
			12	6	6		
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		
2	C	1	Total	C	O	0	0
			12	6	6		
2	I	1	Total	C	O	0	0
			12	6	6		
2	J	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula:  $C_6H_{14}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			7	4	3		
3	D	1	Total	C	O	0	0
			7	4	3		
3	E	1	Total	C	O	0	0
			7	4	3		
3	F	1	Total	C	O	0	0
			7	4	3		
3	F	1	Total	C	O	0	0
			7	4	3		
3	G	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	C	1	Total	C	O	0	0
			7	4	3		
3	C	1	Total	C	O	0	0
			9	6	3		
3	I	1	Total	C	O	0	0
			7	4	3		
3	I	1	Total	C	O	0	0
			7	4	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	J	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total	Cl	0	0
			1	1		
4	E	1	Total	Cl	0	0
			1	1		
4	G	2	Total	Cl	0	0
			2	2		
4	H	1	Total	Cl	0	0
			1	1		
4	A	1	Total	Cl	0	0
			1	1		
4	B	2	Total	Cl	0	0
			2	2		
4	C	1	Total	Cl	0	0
			1	1		
4	I	1	Total	Cl	0	0
			1	1		
4	J	1	Total	Cl	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	56	Total	O	0	0
			56	56		
5	E	63	Total	O	0	0
			63	63		
5	F	67	Total	O	0	0
			67	67		
5	G	64	Total	O	0	0
			64	64		
5	H	57	Total	O	0	0
			57	57		
5	A	67	Total	O	0	0
			67	67		
5	B	73	Total	O	0	0
			73	73		

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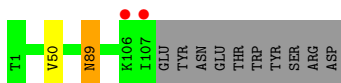
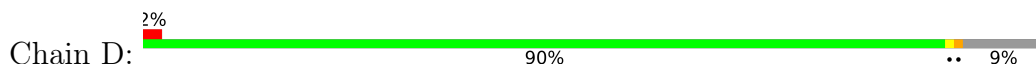
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	66	Total 66	O 66	0	0
5	I	54	Total 54	O 54	0	0
5	J	71	Total 71	O 71	0	0



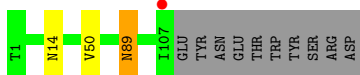
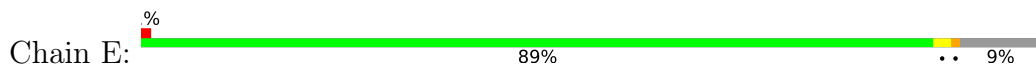
### 3 Residue-property plots [i](#)

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.

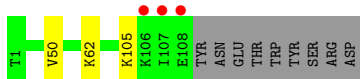
- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



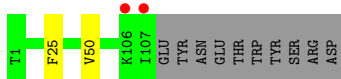
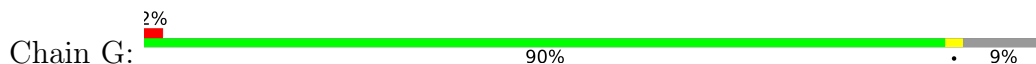
- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



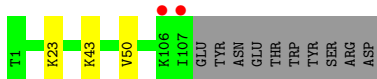
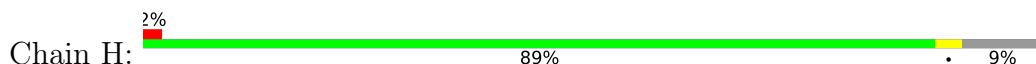
- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



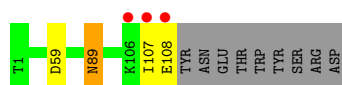
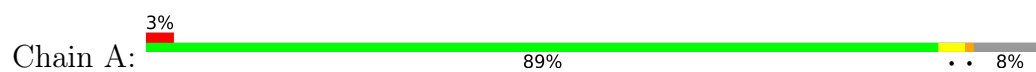
- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



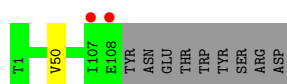
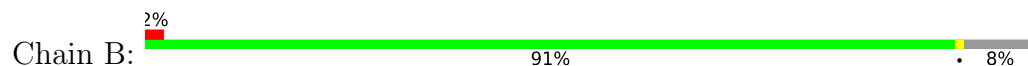
- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



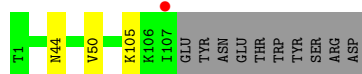
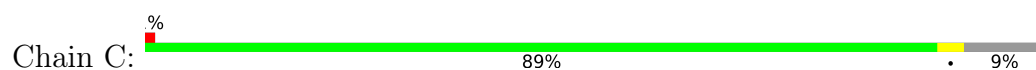
- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



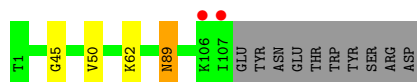
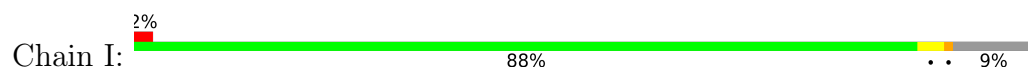
- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



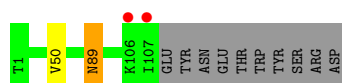
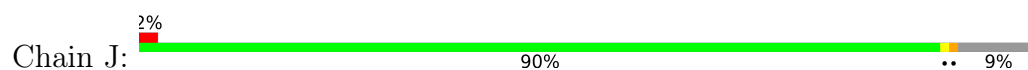
- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



- Molecule 1: Cholera enterotoxin B-subunit, Surface protein gp120



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.29Å 112.43Å 137.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.26 – 2.25 29.26 – 2.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.26-2.25) 100.0 (29.26-2.25)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.26 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, $R_{free}$	0.157 , 0.198 0.164 , 0.201	Depositor DCC
$R_{free}$ test set	1863 reflections (2.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.3	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 29.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	9406	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.87% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, GAL, PGE

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	0.61	0/870	0.66	0/1174
1	B	0.54	0/889	0.62	1/1200 (0.1%)
1	C	0.54	0/878	0.65	1/1184 (0.1%)
1	D	0.48	0/861	0.61	1/1162 (0.1%)
1	E	0.49	0/861	0.61	1/1162 (0.1%)
1	F	0.56	0/870	0.63	1/1174 (0.1%)
1	G	0.48	0/861	0.60	1/1162 (0.1%)
1	H	0.56	0/861	0.61	1/1162 (0.1%)
1	I	0.53	0/872	0.64	1/1177 (0.1%)
1	J	0.51	0/861	0.59	1/1162 (0.1%)
All	All	0.53	0/8684	0.62	9/11719 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	50	VAL	N-CA-C	-6.02	94.75	111.00
1	G	50	VAL	N-CA-C	-5.99	94.84	111.00
1	I	50	VAL	N-CA-C	-5.97	94.87	111.00
1	H	50	VAL	N-CA-C	-5.88	95.12	111.00
1	E	50	VAL	N-CA-C	-5.84	95.22	111.00
1	B	50	VAL	N-CA-C	-5.71	95.57	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	50	VAL	N-CA-C	-5.38	96.47	111.00
1	J	50	VAL	N-CA-C	-5.27	96.78	111.00
1	C	50	VAL	N-CA-C	-5.00	97.50	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	44[A]	ASN	Mainchain

## 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	856	0	866	3	0
1	B	868	0	877	0	0
1	C	861	0	878	1	0
1	D	847	0	860	1	0
1	E	847	0	860	2	0
1	F	856	0	866	2	0
1	G	847	0	860	2	0
1	H	847	0	860	1	0
1	I	854	0	867	2	0
1	J	847	0	860	1	0
2	A	12	0	12	0	0
2	B	12	0	12	0	0
2	C	12	0	12	0	0
2	D	12	0	12	0	0
2	E	12	0	12	0	0
2	F	12	0	12	0	0
2	G	12	0	12	0	0
2	H	12	0	12	0	0
2	I	12	0	12	0	0
2	J	12	0	12	0	0
3	A	14	0	18	0	0
3	B	14	0	18	0	0
3	C	16	0	20	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	14	0	18	0	0
3	E	7	0	9	0	0
3	F	14	0	18	0	0
3	G	7	0	9	0	0
3	I	14	0	18	1	0
3	J	7	0	9	0	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	G	2	0	0	0	0
4	H	1	0	0	0	0
4	I	1	0	0	0	0
4	J	1	0	0	0	0
5	A	67	0	0	1	0
5	B	73	0	0	0	0
5	C	66	0	0	1	0
5	D	56	0	0	0	0
5	E	63	0	0	1	0
5	F	67	0	0	0	0
5	G	64	0	0	0	0
5	H	57	0	0	0	0
5	I	54	0	0	0	0
5	J	71	0	0	0	0
All	All	9406	0	8911	13	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 1.

All (13) 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:107:ILE:O	1:A:108:GLU:HG2	1.64	0.96
1:D:89:ASN:C	1:D:89:ASN:HD22	2.12	0.53
1:C:105[A]:LYS:NZ	5:C:302:HOH:O	2.41	0.52
1:A:89:ASN:ND2	5:A:301:HOH:O	2.40	0.48
1:A:89:ASN:HD22	1:A:89:ASN:C	2.17	0.48
1:J:89:ASN:C	1:J:89:ASN:HD22	2.17	0.47
1:E:89:ASN:C	1:E:89:ASN:HD22	2.18	0.46
1:I:89:ASN:HD22	1:I:89:ASN:C	2.22	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:105:LYS:HE3	1:G:25:PHE:HE2	1.84	0.42
1:E:14:ASN:ND2	5:E:301:HOH:O	2.50	0.42
1:F:105:LYS:HE3	1:G:25:PHE:CE2	2.54	0.42
1:H:43:LYS:HE2	1:H:43:LYS:HB3	1.80	0.42
1:I:45:GLY:HA2	3:I:203:PGE:H4	2.04	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	106/117 (91%)	104 (98%)	2 (2%)	0	100	100
1	B	108/117 (92%)	106 (98%)	2 (2%)	0	100	100
1	C	107/117 (92%)	105 (98%)	2 (2%)	0	100	100
1	D	105/117 (90%)	103 (98%)	2 (2%)	0	100	100
1	E	105/117 (90%)	103 (98%)	2 (2%)	0	100	100
1	F	106/117 (91%)	103 (97%)	3 (3%)	0	100	100
1	G	105/117 (90%)	103 (98%)	2 (2%)	0	100	100
1	H	105/117 (90%)	103 (98%)	2 (2%)	0	100	100
1	I	106/117 (91%)	104 (98%)	2 (2%)	0	100	100
1	J	105/117 (90%)	101 (96%)	4 (4%)	0	100	100
All	All	1058/1170 (90%)	1035 (98%)	23 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	94/103 (91%)	92 (98%)	2 (2%)	53	62
1	B	96/103 (93%)	96 (100%)	0	100	100
1	C	95/103 (92%)	95 (100%)	0	100	100
1	D	93/103 (90%)	92 (99%)	1 (1%)	73	82
1	E	93/103 (90%)	92 (99%)	1 (1%)	73	82
1	F	94/103 (91%)	93 (99%)	1 (1%)	73	82
1	G	93/103 (90%)	93 (100%)	0	100	100
1	H	93/103 (90%)	92 (99%)	1 (1%)	73	82
1	I	94/103 (91%)	92 (98%)	2 (2%)	53	62
1	J	93/103 (90%)	92 (99%)	1 (1%)	73	82
All	All	938/1030 (91%)	929 (99%)	9 (1%)	76	84

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	89	ASN
1	E	89	ASN
1	F	62	LYS
1	H	23	LYS
1	A	59	ASP
1	A	89	ASN
1	I	62	LYS
1	I	89	ASN
1	J	89	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	16	GLN
1	D	89	ASN

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Mol	Chain	Res	Type
1	E	16	GLN
1	E	89	ASN
1	E	103	ASN
1	G	4	ASN
1	G	103	ASN
1	A	16	GLN
1	A	89	ASN
1	C	49	GLN
1	I	4	ASN
1	I	16	GLN
1	I	89	ASN
1	I	103	ASN
1	J	16	GLN
1	J	89	ASN

### 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 36 ligands modelled in this entry, 11 are monoatomic - leaving 25 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
3	PGE	E	202	-	6,6,9	0.53	0	5,5,8	0.50	0
3	PGE	A	202	-	6,6,9	0.44	0	5,5,8	0.30	0
3	PGE	D	202	-	6,6,9	0.42	0	5,5,8	0.31	0
2	GAL	A	201	-	12,12,12	0.49	0	17,17,17	0.90	0
2	GAL	E	201	-	12,12,12	0.63	0	17,17,17	0.79	0
3	PGE	G	202	-	6,6,9	0.45	0	5,5,8	0.30	0
2	GAL	I	201	-	12,12,12	0.69	0	17,17,17	0.61	0
3	PGE	B	203	-	6,6,9	0.41	0	5,5,8	0.35	0
2	GAL	J	201	-	12,12,12	0.67	0	17,17,17	0.70	0
3	PGE	C	203	-	8,8,9	0.53	0	7,7,8	0.18	0
3	PGE	B	202	-	6,6,9	0.50	0	5,5,8	0.33	0
3	PGE	F	203	-	6,6,9	0.65	0	5,5,8	0.54	0
2	GAL	G	201	-	12,12,12	0.61	0	17,17,17	0.73	0
2	GAL	H	201	-	12,12,12	0.53	0	17,17,17	0.72	0
3	PGE	D	203	-	6,6,9	0.46	0	5,5,8	0.32	0
3	PGE	C	202	-	6,6,9	0.37	0	5,5,8	0.34	0
3	PGE	F	202	-	6,6,9	0.46	0	5,5,8	0.32	0
2	GAL	C	201	-	12,12,12	0.56	0	17,17,17	0.78	0
3	PGE	J	202	-	6,6,9	0.44	0	5,5,8	0.24	0
3	PGE	I	202	-	6,6,9	0.46	0	5,5,8	0.39	0
2	GAL	F	201	-	12,12,12	0.62	0	17,17,17	0.73	0
2	GAL	B	201	-	12,12,12	0.53	0	17,17,17	0.80	1 (5%)
3	PGE	A	203	-	6,6,9	0.49	0	5,5,8	0.20	0
3	PGE	I	203	-	6,6,9	0.47	0	5,5,8	0.32	0
2	GAL	D	201	-	12,12,12	0.63	0	17,17,17	0.77	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
3	PGE	E	202	-	-	2/4/4/7	-
3	PGE	A	202	-	-	2/4/4/7	-
3	PGE	D	202	-	-	3/4/4/7	-
2	GAL	A	201	-	-	0/2/22/22	0/1/1/1
2	GAL	E	201	-	-	0/2/22/22	0/1/1/1
3	PGE	G	202	-	-	2/4/4/7	-
2	GAL	I	201	-	-	0/2/22/22	0/1/1/1
3	PGE	B	203	-	-	3/4/4/7	-
2	GAL	J	201	-	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PGE	C	203	-	-	4/6/6/7	-
3	PGE	B	202	-	-	3/4/4/7	-
3	PGE	F	203	-	-	3/4/4/7	-
2	GAL	G	201	-	-	0/2/22/22	0/1/1/1
2	GAL	H	201	-	-	0/2/22/22	0/1/1/1
3	PGE	D	203	-	-	1/4/4/7	-
3	PGE	C	202	-	-	3/4/4/7	-
3	PGE	F	202	-	-	3/4/4/7	-
2	GAL	C	201	-	-	0/2/22/22	0/1/1/1
3	PGE	J	202	-	-	1/4/4/7	-
3	PGE	I	202	-	-	3/4/4/7	-
2	GAL	F	201	-	-	0/2/22/22	0/1/1/1
2	GAL	B	201	-	-	0/2/22/22	0/1/1/1
3	PGE	A	203	-	-	3/4/4/7	-
3	PGE	I	203	-	-	3/4/4/7	-
2	GAL	D	201	-	-	0/2/22/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	GAL	O5-C5-C6	2.06	111.57	106.44

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	203	PGE	O1-C1-C2-O2
3	A	203	PGE	O1-C1-C2-O2
3	B	202	PGE	O1-C1-C2-O2
3	E	202	PGE	C4-C3-O2-C2
3	C	203	PGE	O2-C3-C4-O3
3	A	203	PGE	O2-C3-C4-O3
3	B	203	PGE	O2-C3-C4-O3
3	C	203	PGE	O1-C1-C2-O2
3	I	202	PGE	O1-C1-C2-O2
3	I	203	PGE	O2-C3-C4-O3
3	D	202	PGE	O1-C1-C2-O2
3	G	202	PGE	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	C	202	PGE	O2-C3-C4-O3
3	I	203	PGE	O3-C5-C6-O4
3	F	203	PGE	C1-C2-O2-C3
3	F	202	PGE	C1-C2-O2-C3
3	F	203	PGE	O2-C3-C4-O3
3	B	202	PGE	C4-C3-O2-C2
3	A	202	PGE	O1-C1-C2-O2
3	A	202	PGE	O2-C3-C4-O3
3	B	203	PGE	O1-C1-C2-O2
3	J	202	PGE	O2-C3-C4-O3
3	D	202	PGE	O2-C3-C4-O3
3	C	202	PGE	O1-C1-C2-O2
3	C	203	PGE	C6-C5-O3-C4
3	B	203	PGE	C1-C2-O2-C3
3	C	202	PGE	C1-C2-O2-C3
3	B	202	PGE	O2-C3-C4-O3
3	G	202	PGE	C1-C2-O2-C3
3	I	203	PGE	C3-C4-O3-C5
3	A	203	PGE	C1-C2-O2-C3
3	I	202	PGE	C1-C2-O2-C3
3	I	202	PGE	C4-C3-O2-C2
3	E	202	PGE	O1-C1-C2-O2
3	C	203	PGE	C3-C4-O3-C5
3	F	202	PGE	O2-C3-C4-O3
3	F	202	PGE	O1-C1-C2-O2
3	F	203	PGE	C4-C3-O2-C2
3	D	202	PGE	C1-C2-O2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	203	PGE	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	108/117 (92%)	-0.32	3 (2%) 53 55	25, 33, 50, 105	0
1	B	108/117 (92%)	-0.26	2 (1%) 66 69	26, 35, 48, 110	0
1	C	107/117 (91%)	-0.33	1 (0%) 84 85	27, 34, 49, 92	0
1	D	107/117 (91%)	-0.39	2 (1%) 66 69	27, 35, 50, 84	0
1	E	107/117 (91%)	-0.32	1 (0%) 84 85	26, 35, 48, 85	0
1	F	108/117 (92%)	-0.16	3 (2%) 53 55	26, 33, 51, 110	0
1	G	107/117 (91%)	-0.26	2 (1%) 66 69	26, 35, 50, 85	0
1	H	107/117 (91%)	-0.26	2 (1%) 66 69	28, 39, 49, 88	0
1	I	107/117 (91%)	-0.06	2 (1%) 66 69	28, 37, 51, 93	0
1	J	107/117 (91%)	-0.21	2 (1%) 66 69	26, 33, 47, 76	0
All	All	1073/1170 (91%)	-0.26	20 (1%) 66 69	25, 35, 51, 110	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	107	ILE	9.8
1	F	108	GLU	6.3
1	H	107	ILE	6.0
1	B	108	GLU	5.9
1	F	107	ILE	5.6
1	G	107	ILE	5.0
1	B	107	ILE	5.0
1	J	107	ILE	4.5
1	I	106	LYS	4.2
1	E	107	ILE	4.2
1	D	107	ILE	3.8
1	A	107	ILE	3.8
1	A	108	GLU	3.8

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Mol	Chain	Res	Type	RSRZ
1	J	106	LYS	3.4
1	G	106	LYS	3.2
1	A	106	LYS	2.6
1	F	106	LYS	2.6
1	D	106	LYS	2.5
1	H	106	LYS	2.5
1	C	107	ILE	2.1

## 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
3	PGE	F	203	7/10	0.77	0.29	46,50,54,59	0
3	PGE	A	203	7/10	0.80	0.24	55,64,68,70	0
3	PGE	I	203	7/10	0.80	0.18	68,69,74,75	0
3	PGE	B	203	7/10	0.81	0.28	58,60,62,65	0
3	PGE	C	203	9/10	0.82	0.23	61,67,73,74	0
3	PGE	F	202	7/10	0.86	0.20	57,60,62,62	0
3	PGE	G	202	7/10	0.86	0.15	57,60,66,68	0
3	PGE	J	202	7/10	0.86	0.28	64,66,68,68	0
3	PGE	C	202	7/10	0.87	0.14	53,58,61,62	0
3	PGE	A	202	7/10	0.87	0.17	53,54,61,64	0
3	PGE	I	202	7/10	0.89	0.28	60,62,66,67	0
3	PGE	E	202	7/10	0.91	0.18	52,52,58,59	0
3	PGE	D	203	7/10	0.91	0.20	57,58,63,64	0
3	PGE	B	202	7/10	0.91	0.12	61,62,64,66	0
2	GAL	B	201	12/12	0.93	0.12	30,35,40,43	0
2	GAL	E	201	12/12	0.93	0.12	33,35,39,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GAL	H	201	12/12	0.93	0.09	37,39,43,48	0
4	CL	H	202	1/1	0.93	0.10	62,62,62,62	0
3	PGE	D	202	7/10	0.94	0.14	52,53,56,59	0
2	GAL	A	201	12/12	0.94	0.11	28,29,34,40	0
2	GAL	F	201	12/12	0.94	0.12	28,33,38,39	0
2	GAL	J	201	12/12	0.94	0.10	29,32,37,48	0
2	GAL	C	201	12/12	0.95	0.11	28,31,37,41	0
2	GAL	G	201	12/12	0.95	0.12	30,34,37,41	0
2	GAL	D	201	12/12	0.95	0.12	32,35,39,48	0
4	CL	B	205	1/1	0.95	0.25	64,64,64,64	0
2	GAL	I	201	12/12	0.96	0.11	35,37,44,45	0
4	CL	G	203	1/1	0.96	0.05	61,61,61,61	0
4	CL	E	203	1/1	0.97	0.08	54,54,54,54	0
4	CL	C	204	1/1	0.97	0.19	62,62,62,62	0
4	CL	A	204	1/1	0.98	0.15	54,54,54,54	0
4	CL	B	204	1/1	0.98	0.16	58,58,58,58	0
4	CL	G	204	1/1	0.98	0.05	43,43,43,43	0
4	CL	D	204	1/1	0.98	0.21	58,58,58,58	0
4	CL	I	204	1/1	0.98	0.21	58,58,58,58	0
4	CL	J	203	1/1	0.98	0.11	61,61,61,61	0

## 6.5 Other polymers

There are no such residues in this entry.