



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 12, 2017 – 08:46 AM EST

PDB ID : 5TPB  
Title : Binding domain of BoNT/A complexed with ganglioside variant  
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Deposited on : 2016-10-20  
Resolution : 2.60 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20028442

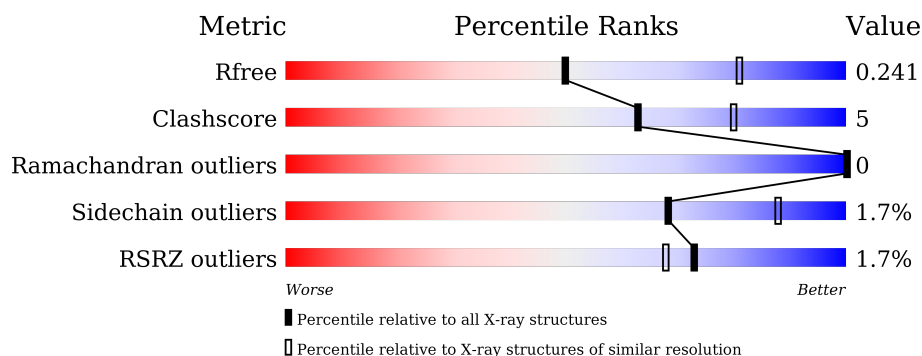
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	444	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>14%</div> <div>7%</div> </div> </div>
1	B	444	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>12%</div> <div>8%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SIA	B	2301	X	-	-	X
3	GAL	B	2302	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6903 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 Botulinum neurotoxin type A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	0	0
			3371	2148	576	633	14			
1	B	407	Total	C	N	O	S	0	2	0
			3375	2154	581	627	13			

There are 46 discrepancies between the modelled and reference sequences:

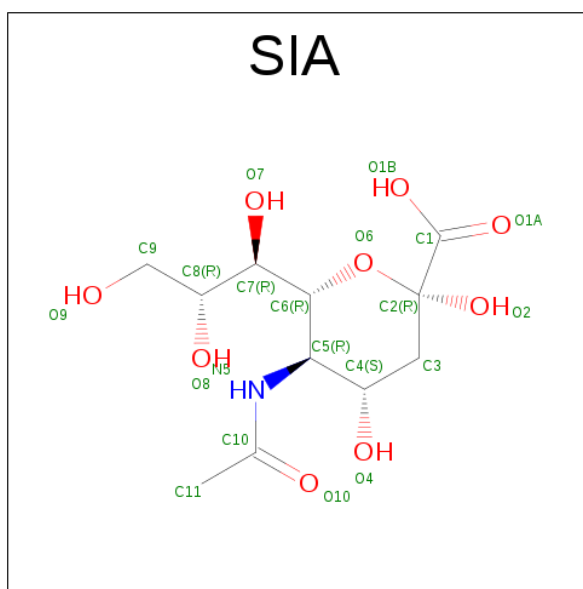
Chain	Residue	Modelled	Actual	Comment	Reference
A	854	MET	-	initiating methionine	UNP P10845
A	855	GLY	-	expression tag	UNP P10845
A	856	SER	-	expression tag	UNP P10845
A	857	SER	-	expression tag	UNP P10845
A	858	HIS	-	expression tag	UNP P10845
A	859	HIS	-	expression tag	UNP P10845
A	860	HIS	-	expression tag	UNP P10845
A	861	HIS	-	expression tag	UNP P10845
A	862	HIS	-	expression tag	UNP P10845
A	863	HIS	-	expression tag	UNP P10845
A	864	SER	-	expression tag	UNP P10845
A	865	SER	-	expression tag	UNP P10845
A	866	GLY	-	expression tag	UNP P10845
A	867	LEU	-	expression tag	UNP P10845
A	868	VAL	-	expression tag	UNP P10845
A	869	PRO	-	expression tag	UNP P10845
A	870	ARG	-	expression tag	UNP P10845
A	871	GLY	-	expression tag	UNP P10845
A	872	SER	-	expression tag	UNP P10845
A	873	HIS	-	expression tag	UNP P10845
A	874	MET	-	expression tag	UNP P10845
A	875	ASP	-	expression tag	UNP P10845
A	1297	GLN	-	expression tag	UNP P10845
B	854	MET	-	initiating methionine	UNP P10845
B	855	GLY	-	expression tag	UNP P10845

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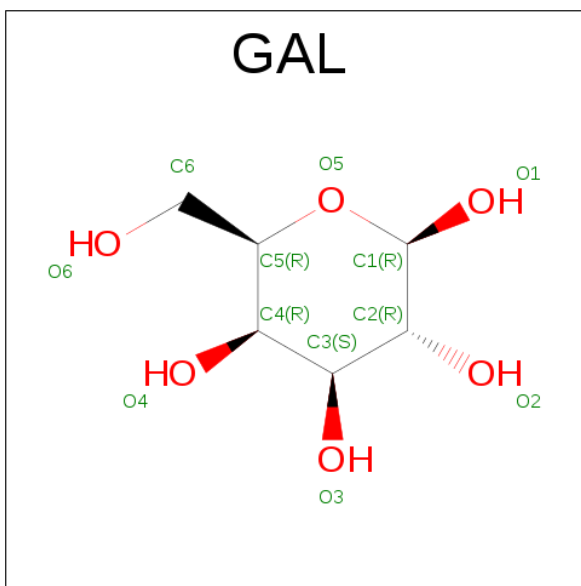
Chain	Residue	Modelled	Actual	Comment	Reference
B	856	SER	-	expression tag	UNP P10845
B	857	SER	-	expression tag	UNP P10845
B	858	HIS	-	expression tag	UNP P10845
B	859	HIS	-	expression tag	UNP P10845
B	860	HIS	-	expression tag	UNP P10845
B	861	HIS	-	expression tag	UNP P10845
B	862	HIS	-	expression tag	UNP P10845
B	863	HIS	-	expression tag	UNP P10845
B	864	SER	-	expression tag	UNP P10845
B	865	SER	-	expression tag	UNP P10845
B	866	GLY	-	expression tag	UNP P10845
B	867	LEU	-	expression tag	UNP P10845
B	868	VAL	-	expression tag	UNP P10845
B	869	PRO	-	expression tag	UNP P10845
B	870	ARG	-	expression tag	UNP P10845
B	871	GLY	-	expression tag	UNP P10845
B	872	SER	-	expression tag	UNP P10845
B	873	HIS	-	expression tag	UNP P10845
B	874	MET	-	expression tag	UNP P10845
B	875	ASP	-	expression tag	UNP P10845
B	1297	GLN	-	expression tag	UNP P10845

- Molecule 2 is O-SIALIC ACID (three-letter code: SIA) (formula: C<sub>11</sub>H<sub>19</sub>NO<sub>9</sub>).



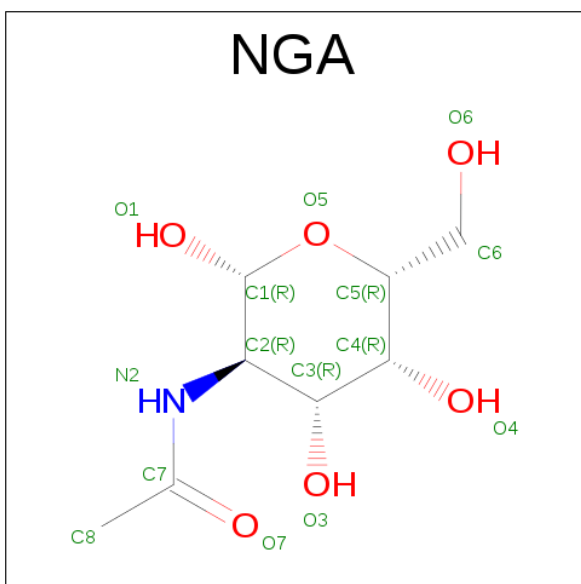
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			20	11	1	8		

- Molecule 3 is BETA-D-GALACTOSE (three-letter code: GAL) (formula:  $C_6H_{12}O_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			11	6	5		

- Molecule 4 is N-ACETYL-D-GALACTOSAMINE (three-letter code: NGA) (formula:  $C_8H_{15}NO_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			15	8	1	6		

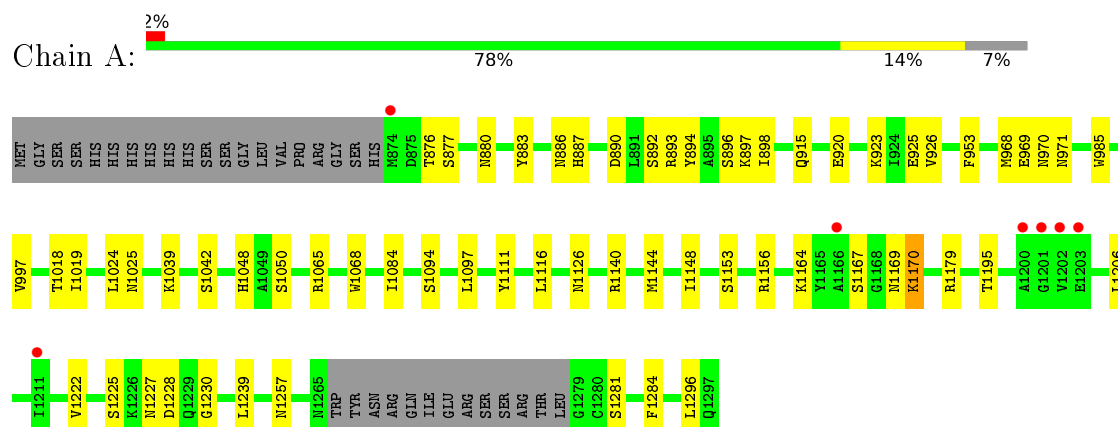
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	58	Total 58	O 58	0	0
5	B	53	Total 53	O 53	0	0

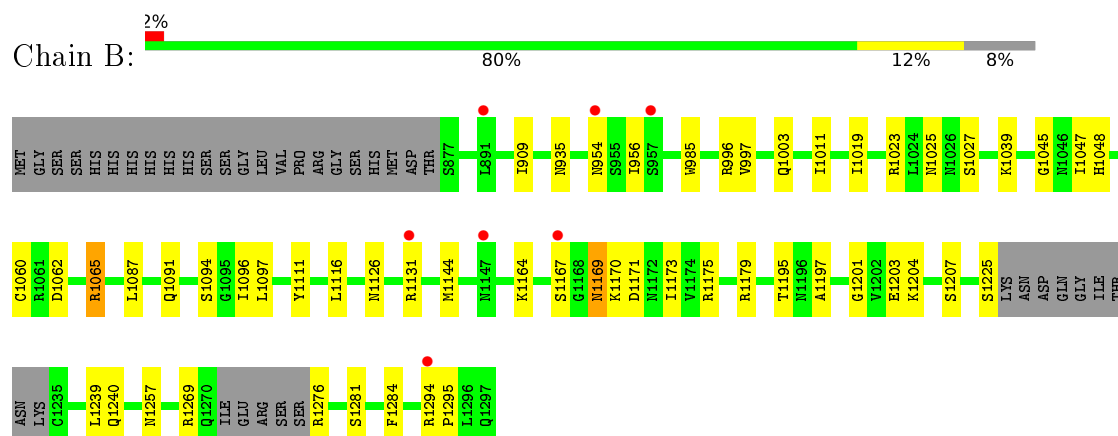
### 3 Residue-property plots [i](#)

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

#### • Molecule 1: Botulinum neurotoxin type A



#### • Molecule 1: Botulinum neurotoxin type A





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.52Å 104.33Å 69.41Å 90.00° 115.97° 90.00°	Depositor
Resolution (Å)	39.05 – 2.60 39.05 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (39.05-2.60) 99.7 (39.05-2.60)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.73 (at 2.61Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.222 , 0.240 0.225 , 0.241	Depositor DCC
$R_{free}$ test set	1325 reflections (5.41%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.7	Xtriage
Anisotropy	0.630	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 34.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.034 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	6903	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.55% 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: SIA, GAL, NGA

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	0.56	0/3440	0.58	0/4648
1	B	0.51	1/3446 (0.0%)	0.59	1/4656 (0.0%)
All	All	0.54	1/6886 (0.0%)	0.58	1/9304 (0.0%)

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	A	0	8
1	B	0	3
All	All	0	11

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1295	PRO	N-CD	5.06	1.54	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1294	ARG	C-N-CD	5.79	140.56	128.40

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1094	SER	Peptide

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Mol	Chain	Res	Type	Group
1	A	1144	MET	Peptide
1	A	1167	SER	Peptide
1	A	1170	LYS	Peptide
1	A	1227	ASN	Peptide
1	A	1228	ASP	Peptide
1	A	1230	GLY	Peptide
1	A	1257	ASN	Peptide
1	B	1144	MET	Peptide
1	B	1257	ASN	Peptide
1	B	996	ARG	Peptide

## 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	3371	0	3329	33	0
1	B	3375	0	3330	32	0
2	B	20	0	17	7	0
3	B	11	0	10	8	0
4	B	15	0	14	0	0
5	A	58	0	0	1	0
5	B	53	0	0	3	0
All	All	6903	0	6700	71	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 (71) 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)
2:B:2301:SIA:C2	3:B:2302:GAL:O3	1.71	1.38
1:B:1269:ARG:HG3	5:B:2443:HOH:O	1.01	1.18
2:B:2301:SIA:C3	3:B:2302:GAL:O3	2.06	1.01
1:A:892:SER:OG	1:A:894:TYR:CD2	2.27	0.87
1:B:1126:ASN:HB3	5:B:2437:HOH:O	1.77	0.84
2:B:2301:SIA:H32	3:B:2302:GAL:O3	1.76	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:2301:SIA:C2	3:B:2302:GAL:HO3	1.90	0.83
1:A:1140:ARG:HD3	1:A:1153:SER:O	1.77	0.82
1:A:969:GLU:HG3	1:A:1050:SER:HA	1.61	0.82
1:B:1197:ALA:O	1:B:1204:LYS:NZ	2.13	0.81
1:B:1164:LYS:NZ	1:B:1169:ASN:O	2.15	0.80
1:B:1096:ILE:HD13	1:B:1173:ILE:HG21	1.61	0.80
1:B:1060:CYS:SG	1:B:1065:ARG:HG3	2.27	0.73
1:B:935:ASN:HB2	1:B:1048:HIS:CE1	2.24	0.73
1:B:1179:ARG:NH2	1:B:1240:GLN:OE1	2.24	0.71
2:B:2301:SIA:C1	3:B:2302:GAL:O3	2.39	0.68
1:B:1203:GLU:OE2	3:B:2302:GAL:O6	2.11	0.67
1:B:909:ILE:HG23	1:B:1173:ILE:HD11	1.77	0.67
1:B:1065:ARG:O	5:B:2401:HOH:O	2.13	0.65
1:A:1164:LYS:NZ	1:A:1169:ASN:O	2.30	0.64
1:B:1201:GLY:O	1:B:1204:LYS:HE2	2.01	0.60
1:B:935:ASN:HB2	1:B:1048:HIS:ND1	2.16	0.60
2:B:2301:SIA:C1	3:B:2302:GAL:HO3	2.13	0.59
1:B:1097:LEU:HD22	1:B:1225:SER:HB3	1.85	0.59
1:A:925:GLU:OE1	1:B:1276:ARG:N	2.36	0.58
1:A:892:SER:OG	1:A:894:TYR:HD2	1.85	0.58
1:A:970:ASN:O	1:A:971:ASN:HB2	2.03	0.58
1:A:969:GLU:HG2	1:A:1050:SER:HB2	1.84	0.58
1:B:1023:ARG:NH2	1:B:1045:GLY:O	2.34	0.57
1:A:968:MET:O	1:A:969:GLU:HG2	2.05	0.56
1:A:898:ILE:HD12	1:A:926:VAL:HG22	1.87	0.55
1:A:1024:LEU:O	1:A:1042:SER:HB3	2.07	0.55
1:A:1179:ARG:HG2	1:A:1222:VAL:HG22	1.89	0.54
1:A:876:THR:HG22	1:A:877:SER:N	2.21	0.54
1:A:896:SER:O	1:A:897:LYS:HB3	2.08	0.54
1:A:1097:LEU:HD22	1:A:1225:SER:HB3	1.90	0.53
1:B:1167:SER:OG	1:B:1169:ASN:HB2	2.09	0.52
1:A:1097:LEU:CD2	1:A:1225:SER:HB3	2.42	0.50
1:A:1116:LEU:HD12	1:A:1281:SER:HB3	1.94	0.50
1:B:1087:LEU:O	1:B:1091:GLN:HG3	2.12	0.50
1:A:883:TYR:CZ	1:A:886:ASN:HA	2.48	0.49
1:A:1195:THR:HB	1:A:1206:LEU:HD23	1.95	0.48
1:B:1131:ARG:HG2	1:B:1131:ARG:O	2.13	0.48
1:A:890:ASP:OD1	1:A:892:SER:HB3	2.13	0.48
1:B:1062:ASP:HB3	1:B:1065:ARG:HG2	1.97	0.47
1:A:877:SER:OG	1:A:880:ASN:HB2	2.16	0.46
1:B:1116:LEU:HD12	1:B:1281:SER:HB3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1171:ASP:OD2	1:B:1175:ARG:NH2	2.41	0.45
1:B:1171:ASP:OD1	1:B:1171:ASP:N	2.49	0.45
1:A:883:TYR:CE1	1:A:886:ASN:HA	2.51	0.44
1:A:893:ARG:O	1:A:894:TYR:CD1	2.70	0.44
1:B:997:VAL:HG12	1:B:1039:LYS:HG3	2.00	0.44
1:B:985:TRP:CD2	1:B:1019:ILE:HG21	2.52	0.44
1:A:1140:ARG:HB2	1:A:1156:ARG:HG3	1.99	0.44
1:A:887:HIS:CD2	5:A:1355:HOH:O	2.71	0.44
1:A:997:VAL:HG12	1:A:1039:LYS:HG3	1.99	0.44
1:B:954:ASN:HB3	1:B:956:ILE:HG22	1.98	0.44
2:B:2301:SIA:C3	3:B:2302:GAL:HO3	2.18	0.43
1:B:1111:TYR:CG	1:B:1284:PHE:HB3	2.54	0.43
1:A:985:TRP:CD2	1:A:1019:ILE:HG21	2.53	0.43
1:B:1195:THR:HG22	1:B:1239:LEU:HD21	2.01	0.43
1:B:1023:ARG:NH1	1:B:1047:ILE:HD12	2.33	0.42
1:A:915:GLN:HB2	1:A:1068:TRP:CZ3	2.53	0.42
1:A:1195:THR:HG22	1:A:1239:LEU:HD21	2.01	0.42
1:A:1111:TYR:CG	1:A:1284:PHE:HB3	2.55	0.41
1:A:953:PHE:O	1:A:1148:ILE:HG21	2.20	0.41
1:B:1276:ARG:HD2	1:B:1276:ARG:HA	1.92	0.41
1:A:920:GLU:O	1:A:923:LYS:HD3	2.20	0.41
1:A:1018:THR:HG21	1:A:1084:ILE:HG12	2.03	0.41
1:B:1096:ILE:HG23	1:B:1173:ILE:CG2	2.51	0.40
1:B:1003:GLN:HA	1:B:1011:ILE:HD11	2.03	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	407/444 (92%)	382 (94%)	25 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	403/444 (91%)	385 (96%)	18 (4%)	0	100	100
All	All	810/888 (91%)	767 (95%)	43 (5%)	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	378/408 (93%)	372 (98%)	6 (2%)	70	89
1	B	377/408 (92%)	370 (98%)	7 (2%)	65	86
All	All	755/816 (92%)	742 (98%)	13 (2%)	68	88

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1025	ASN
1	A	1048	HIS
1	A	1065	ARG
1	A	1126	ASN
1	A	1170	LYS
1	A	1296	LEU
1	B	1025	ASN
1	B	1027	SER
1	B	1065	ARG
1	B	1094	SER
1	B	1169	ASN
1	B	1170	LYS
1	B	1207	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	887	HIS
1	A	988	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

3 ligands are modelled in this entry.

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	SIA	B	2301	-	17,20,21	0.56	0	18,28,31	0.97	2 (11%)
3	GAL	B	2302	4	11,11,12	1.06	1 (9%)	15,15,17	1.89	6 (40%)
4	NGA	B	2303	3	15,15,15	1.00	0	17,21,21	1.97	6 (35%)

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	SIA	B	2301	-	1/1/8/9	0/14/34/38	0/1/1/1
3	GAL	B	2302	4	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NGA	B	2303	3	-	0/6/26/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2302	GAL	O5-C1	-2.38	1.39	1.43

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	2303	NGA	C3-C4-C5	-3.60	103.80	110.23
4	B	2303	NGA	O3-C3-C2	-3.46	102.41	109.57
3	B	2302	GAL	O5-C5-C4	-2.80	105.49	110.13
4	B	2303	NGA	O6-C6-C5	-2.77	102.06	111.30
4	B	2303	NGA	O5-C5-C4	-2.75	104.42	109.67
4	B	2303	NGA	C3-C2-N2	-2.71	105.05	110.67
4	B	2303	NGA	O1-C1-O5	-2.67	102.88	110.33
3	B	2302	GAL	O4-C4-C3	-2.65	104.38	110.36
2	B	2301	SIA	C3-C4-C5	-2.46	108.73	111.47
3	B	2302	GAL	O2-C2-C3	-2.33	105.50	110.19
2	B	2301	SIA	O6-C6-C5	-2.24	104.82	108.48
3	B	2302	GAL	C1-O5-C5	2.46	115.76	112.14
3	B	2302	GAL	C1-C2-C3	2.69	112.81	109.55
3	B	2302	GAL	O3-C3-C4	3.01	117.15	110.36

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	2301	SIA	C2

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2301	SIA	7	0
3	B	2302	GAL	8	0



## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	411/444 (92%)	-0.16	7 (1%) 73 68	28, 49, 85, 132	4 (0%)
1	B	407/444 (91%)	-0.10	7 (1%) 73 68	34, 54, 82, 114	4 (0%)
All	All	818/888 (92%)	-0.13	14 (1%) 73 68	28, 52, 84, 132	8 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1202	VAL	5.3
1	A	874	MET	4.2
1	B	957	SER	4.0
1	A	1201	GLY	3.9
1	B	1131	ARG	3.8
1	B	891	LEU	3.1
1	B	1294	ARG	2.8
1	A	1203	GLU	2.8
1	B	954	ASN	2.6
1	B	1167	SER	2.5
1	A	1211	ILE	2.4
1	A	1166	ALA	2.2
1	A	1200	ALA	2.2
1	B	1147	ASN	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	SIA	B	2301	20/21	0.81	0.26	4.59	77,93,102,104	0
3	GAL	B	2302	11/12	0.91	0.14	-0.39	63,68,71,74	0
4	NGA	B	2303	15/15	0.92	0.21	-	71,79,84,90	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.