



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:24 AM GMT

PDB ID : 2WYI  
Title : STRUCTURE OF THE STREPTOCOCCUS PYOGENES FAMILY GH38  
ALPHA-MANNOSIDASE COMPLEXED WITH SWAINSONINE  
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Deposited on : 2009-11-16  
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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

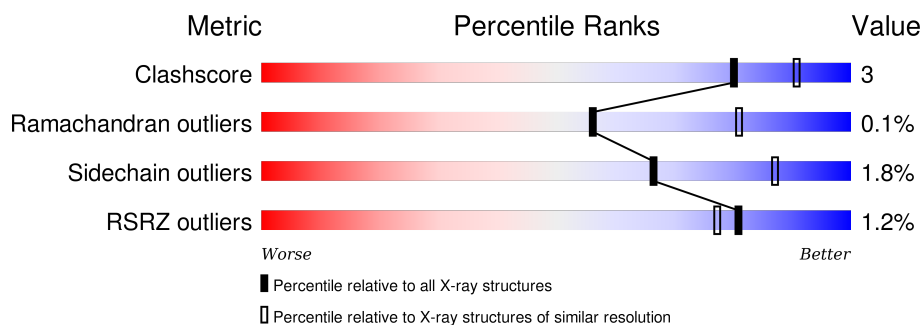
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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))
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	923	<div> <div></div> <div>90%</div> <div>7%</div> <div></div> </div>
1	B	923	<div> <div></div> <div>88%</div> <div>8%</div> <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	PG0	A	1902	-	-	-	X
2	PG0	A	1904	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PG0	A	1905	-	-	-	X
2	PG0	B	1902	-	-	-	X
2	PG0	B	1903	-	-	-	X
2	PG0	B	1904	-	-	-	X
2	PG0	B	1906	-	-	-	X
4	SWA	A	1907	-	-	-	X
4	SWA	B	1909	-	-	-	X

## 2 Entry composition [i](#)

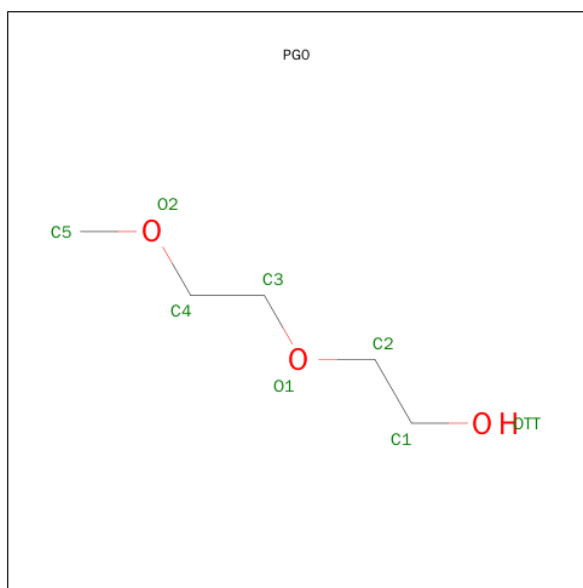
There are 5 unique types of molecules in this entry. The entry contains 14790 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ALPHA-MANNOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	896	Total	C	N	O	S	0	6	0
			7194	4541	1232	1391	30			
1	B	896	Total	C	N	O	S	0	4	0
			7137	4511	1225	1375	26			

- Molecule 2 is 2-(2-METHOXYETHOXY)ETHANOL (three-letter code: PG0) (formula: C<sub>5</sub>H<sub>12</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	5	3		
2	A	1	Total	C	O	0	0
			8	5	3		
2	A	1	Total	C	O	0	0
			8	5	3		
2	A	1	Total	C	O	0	0
			8	5	3		

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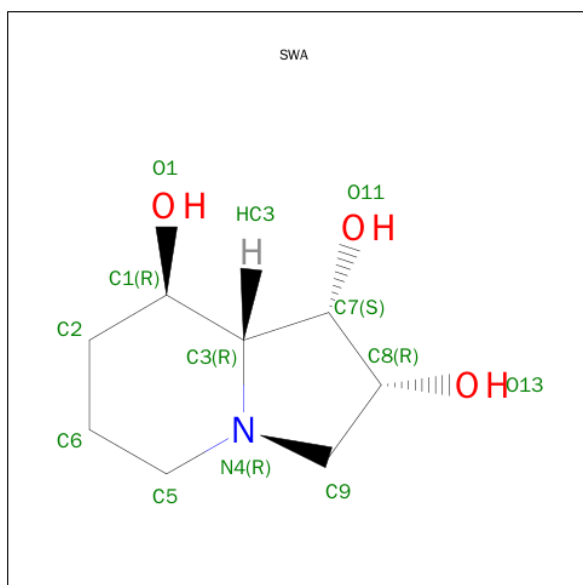
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			8	5	3		
2	B	1	Total	C	O	0	0
			8	5	3		
2	B	1	Total	C	O	0	0
			8	5	3		
2	B	1	Total	C	O	0	0
			8	5	3		
2	B	1	Total	C	O	0	0
			8	5	3		
2	B	1	Total	C	O	0	0
			8	5	3		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is 1S-8AB-OCTAHYDRO-INDOLIZIDINE-1A,2A,8B-TRIOI (three-letter code: SWA) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			12	8	1	3		
4	B	1	Total	C	N	O	0	0
			12	8	1	3		

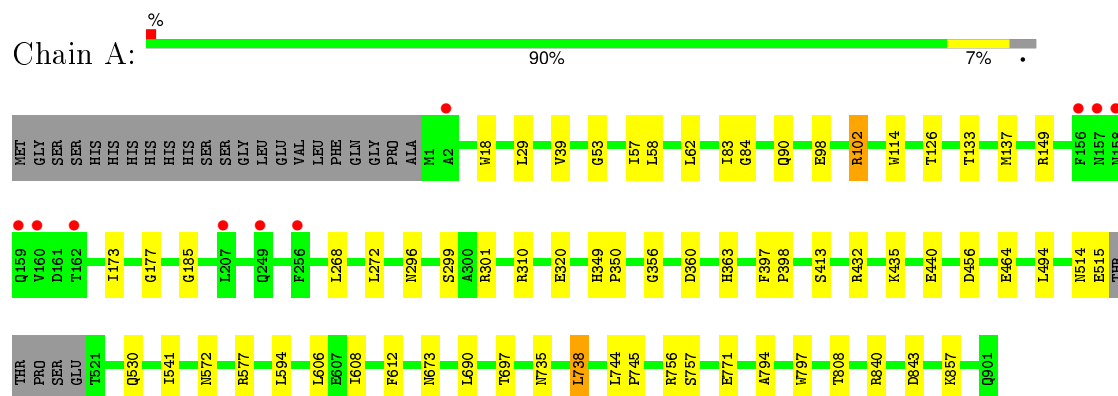
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	222	Total	O	0	0
			222	222		
5	B	123	Total	O	0	0
			123	123		

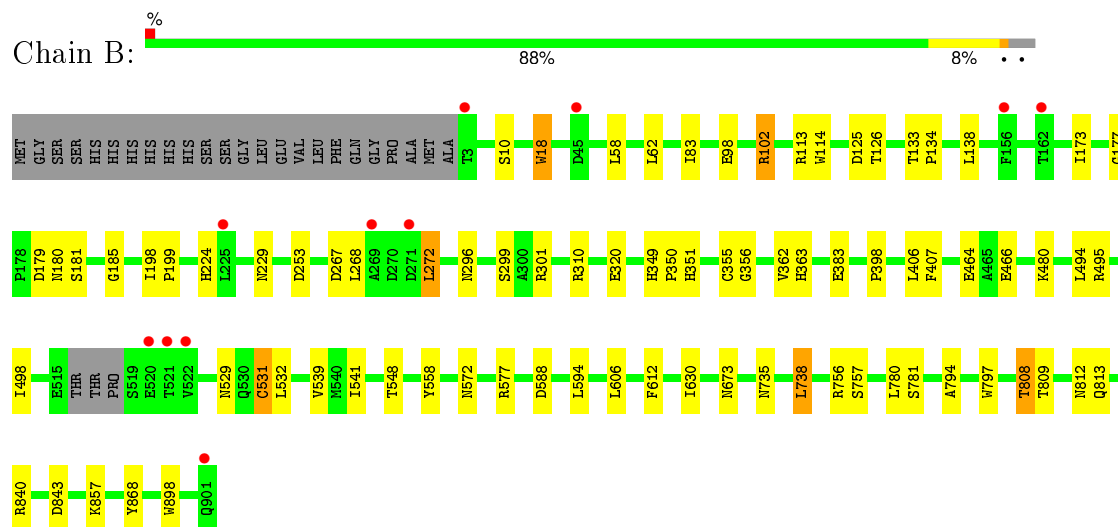
### 3 Residue-property plots [i](#)

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: ALPHA-MANNOSIDASE



#### • Molecule 1: ALPHA-MANNOSIDASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	178.68Å 178.68Å 198.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	132.45 – 2.60 49.56 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (132.45-2.60) 92.9 (49.56-2.60)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.24 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.187 , 0.223 0.185 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	41.7	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 35.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 97828 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	14790	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SWA, ZN, PG0

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	0.47	0/7378	0.59	1/10018 (0.0%)
1	B	0.44	0/7317	0.56	1/9947 (0.0%)
All	All	0.46	0/14695	0.57	2/19965 (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	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	301	ARG	NE-CZ-NH2	-5.51	117.55	120.30
1	A	301	ARG	NE-CZ-NH2	-5.10	117.75	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	90	GLN	Peptide

### 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7194	0	6934	42	0
1	B	7137	0	6824	54	0
2	A	32	0	48	1	0
2	B	56	0	84	5	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	12	0	15	1	0
4	B	12	0	15	3	0
5	A	222	0	0	2	0
5	B	123	0	0	5	0
All	All	14790	0	13920	97	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 3.

All (97) 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:B:812[B]:ASN:HD22	1:B:812[B]:ASN:H	1.16	0.90
1:A:735:ASN:HD22	1:A:738:LEU:H	1.20	0.87
1:A:541:ILE:HD11	1:A:606:LEU:HD23	1.62	0.82
2:B:1902:PG0:H52	5:B:2116:HOH:O	1.84	0.77
1:B:310:ARG:HE	2:B:1903:PG0:H31	1.51	0.74
1:B:529:ASN:HB3	1:B:531:CYS:HB2	1.70	0.72
1:A:735:ASN:ND2	1:A:738:LEU:H	1.87	0.72
1:B:572:ASN:O	1:B:756:ARG:NH2	2.23	0.72
1:B:98:GLU:OE1	1:B:102:ARG:NH1	2.23	0.71
1:B:98:GLU:O	1:B:102:ARG:HG2	1.90	0.70
1:A:98:GLU:O	1:A:102:ARG:HG2	1.91	0.70
1:B:541:ILE:HD11	1:B:606:LEU:HD23	1.72	0.69
1:A:53:GLY:HA2	1:A:84:GLY:O	1.96	0.65
1:B:356:GLY:HA2	1:B:363:HIS:CE1	2.31	0.65
1:B:735:ASN:HB2	1:B:738:LEU:HB2	1.79	0.65
1:B:812[B]:ASN:ND2	1:B:812[B]:ASN:H	1.94	0.62
1:B:398:PRO:O	1:B:813:GLN:NE2	2.32	0.62
1:A:299:SER:OG	1:A:756:ARG:HD3	1.99	0.61
1:A:735:ASN:HB2	1:A:738:LEU:HB2	1.81	0.61
1:A:102:ARG:NH2	1:A:320:GLU:OE2	2.33	0.61
1:A:268:LEU:O	1:A:272:LEU:HB2	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:ASP:OD1	4:B:1909:SWA:HC91	2.01	0.59
1:B:808:THR:HB	5:B:2120:HOH:O	2.02	0.58
1:B:735:ASN:HD22	1:B:738:LEU:H	1.51	0.58
1:A:541:ILE:HD11	1:A:606:LEU:CD2	2.32	0.57
1:B:532:LEU:HB2	1:B:539:VAL:HB	1.86	0.57
1:A:349:HIS:N	1:A:350:PRO:CD	2.68	0.57
1:A:572:ASN:O	1:A:756:ARG:NH2	2.39	0.56
1:B:541:ILE:HD12	1:B:606:LEU:HD21	1.88	0.56
1:B:735:ASN:ND2	1:B:738:LEU:H	2.04	0.55
1:B:781:SER:OG	2:B:1905:PG0:H53	2.07	0.55
1:A:177:GLY:HA3	5:A:2031:HOH:O	2.06	0.55
1:B:541:ILE:HD11	1:B:606:LEU:CD2	2.37	0.55
1:B:466:GLU:OE1	1:B:495:ARG:HD2	2.08	0.54
1:B:383:GLU:CD	1:B:480:LYS:HZ1	2.11	0.54
1:A:356:GLY:HA2	1:A:363:HIS:CE1	2.43	0.54
1:B:268:LEU:HB3	1:B:272:LEU:CD2	2.39	0.53
1:A:173:ILE:HG13	1:A:185:GLY:HA3	1.91	0.53
1:A:541:ILE:CD1	1:A:606:LEU:HD23	2.37	0.53
1:B:102:ARG:NH2	1:B:320:GLU:OE1	2.42	0.53
1:B:840:ARG:HD2	1:B:843:ASP:OD1	2.08	0.53
1:A:296:ASN:O	1:A:756:ARG:NH1	2.43	0.52
1:B:125:ASP:OD1	4:B:1909:SWA:HC52	2.10	0.52
1:B:541:ILE:CD1	1:B:606:LEU:CD2	2.89	0.51
1:A:541:ILE:CD1	1:A:606:LEU:CD2	2.89	0.51
1:A:62:LEU:HD11	1:A:114:TRP:HH2	1.76	0.51
1:B:310:ARG:NE	2:B:1903:PG0:H22	2.26	0.50
1:B:179:ASP:O	1:B:180:ASN:HB2	2.11	0.50
1:B:673:ASN:HA	1:B:757:SER:O	2.11	0.50
1:A:514:ASN:O	1:A:515:GLU:HB2	2.12	0.50
1:A:515:GLU:OE1	1:A:515:GLU:HA	2.11	0.49
1:B:62:LEU:HD11	1:B:114:TRP:HH2	1.77	0.49
1:A:58:LEU:HD11	1:A:83:ILE:HG21	1.95	0.49
1:B:58:LEU:HD11	1:B:83:ILE:HG21	1.94	0.49
1:A:673:ASN:HA	1:A:757:SER:O	2.13	0.48
1:A:149:ARG:NH2	4:A:1907:SWA:HC62	2.27	0.48
1:A:435:LYS:HB3	1:B:630:ILE:HD13	1.95	0.48
1:B:18:TRP:CZ2	4:B:1909:SWA:HC3	2.49	0.47
1:A:29:LEU:HD11	1:A:57:ILE:HG23	1.96	0.47
1:B:299:SER:OG	1:B:756:ARG:HD3	2.15	0.47
1:A:133:THR:O	1:A:137:MET:HG2	2.14	0.47
1:B:351:HIS:O	1:B:355:CYS:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:ILE:HG13	1:B:185:GLY:HA3	1.96	0.46
1:A:794:ALA:HA	1:A:797:TRP:NE1	2.31	0.46
1:B:362:VAL:HA	5:B:2054:HOH:O	2.16	0.46
1:B:794:ALA:HA	1:B:797:TRP:NE1	2.30	0.46
1:A:432:ARG:NH1	1:A:440:GLU:OE1	2.46	0.46
1:A:840:ARG:HD2	1:A:843:ASP:OD1	2.16	0.46
1:B:541:ILE:CD1	1:B:606:LEU:HD21	2.46	0.46
1:A:360:ASP:CG	1:A:771:GLU:HB2	2.35	0.45
1:B:296:ASN:O	1:B:756:ARG:NH1	2.49	0.45
1:B:62:LEU:HD11	1:B:114:TRP:CH2	2.51	0.45
1:B:349:HIS:N	1:B:350:PRO:CD	2.80	0.45
1:B:224:HIS:CE1	1:B:267:ASP:HB3	2.51	0.45
1:A:690:LEU:HB3	1:A:697:THR:HG22	2.00	0.44
1:B:138:LEU:HD12	1:B:181:SER:HB3	1.98	0.44
1:A:102:ARG:HH22	1:A:320:GLU:CD	2.20	0.44
1:A:744:LEU:HA	1:A:745:PRO:HD3	1.77	0.43
1:B:407:PHE:HE2	1:B:809:THR:HG22	1.83	0.43
1:B:840:ARG:CD	1:B:843:ASP:OD1	2.67	0.43
1:A:541:ILE:HD12	1:A:606:LEU:HD21	2.01	0.43
1:B:406:LEU:HD12	1:B:498:ILE:HD13	2.01	0.43
1:A:456:ASP:HB2	5:A:2092:HOH:O	2.19	0.42
1:B:548:THR:HA	1:B:558:TYR:O	2.19	0.42
1:B:177:GLY:HA3	5:B:2030:HOH:O	2.19	0.42
1:A:541:ILE:HD13	1:A:608:ILE:HD11	2.02	0.41
2:B:1902:PG0:H32	5:B:2114:HOH:O	2.20	0.41
1:A:397:PHE:HA	1:A:398:PRO:HD3	1.92	0.41
1:B:868:TYR:HB3	1:B:898:TRP:HB3	2.02	0.41
1:B:198:ILE:HA	1:B:199:PRO:HD3	1.84	0.41
1:A:310:ARG:HE	2:A:1905:PG0:H41	1.84	0.41
1:A:173:ILE:CG1	1:A:185:GLY:HA3	2.51	0.41
1:B:10:SER:HA	1:B:229:ASN:HB3	2.03	0.41
1:A:794:ALA:HA	1:A:797:TRP:CD1	2.56	0.40
1:A:126:THR:O	1:A:149:ARG:HD2	2.21	0.40
1:B:133:THR:HB	1:B:134:PRO:HD3	2.03	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	898/923 (97%)	875 (97%)	22 (2%)	1 (0%)	56	81
1	B	896/923 (97%)	869 (97%)	26 (3%)	1 (0%)	56	81
All	All	1794/1846 (97%)	1744 (97%)	48 (3%)	2 (0%)	56	81

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	TRP
1	B	18	TRP

### 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	780/810 (96%)	768 (98%)	12 (2%)	72	90
1	B	764/810 (94%)	748 (98%)	16 (2%)	61	85
All	All	1544/1620 (95%)	1516 (98%)	28 (2%)	66	87

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	VAL
1	A	102	ARG
1	A	413	SER
1	A	464	GLU

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Mol	Chain	Res	Type
1	A	494	LEU
1	A	530	GLN
1	A	577	ARG
1	A	594	LEU
1	A	612	PHE
1	A	738	LEU
1	A	808	THR
1	A	857	LYS
1	B	102	ARG
1	B	113	ARG
1	B	126	THR
1	B	253	ASP
1	B	272	LEU
1	B	464	GLU
1	B	494	LEU
1	B	531	CYS
1	B	577	ARG
1	B	588	ASP
1	B	594	LEU
1	B	612	PHE
1	B	738	LEU
1	B	780	LEU
1	B	808	THR
1	B	857	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	402	GLN
1	A	453	GLN
1	A	735	ASN
1	B	275	ASN
1	B	530	GLN
1	B	718	GLN
1	B	729	ASN
1	B	735	ASN

### 5.3.3 RNA ⓘ

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 2 are monoatomic - leaving 13 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	PG0	A	1902	-	7,7,7	0.44	0	6,6,6	0.55	0
2	PG0	A	1903	-	7,7,7	0.37	0	6,6,6	0.52	0
2	PG0	A	1904	-	7,7,7	0.62	0	6,6,6	0.29	0
2	PG0	A	1905	-	7,7,7	0.56	0	6,6,6	0.19	0
4	SWA	A	1907	3	13,13,13	0.86	0	13,19,19	1.41	1 (7%)
2	PG0	B	1901	-	7,7,7	0.55	0	6,6,6	0.19	0
2	PG0	B	1902	-	7,7,7	0.52	0	6,6,6	0.10	0
2	PG0	B	1903	-	7,7,7	0.45	0	6,6,6	0.26	0
2	PG0	B	1904	-	7,7,7	0.53	0	6,6,6	0.22	0
2	PG0	B	1905	-	7,7,7	0.48	0	6,6,6	0.31	0
2	PG0	B	1906	-	7,7,7	0.61	0	6,6,6	0.42	0
2	PG0	B	1907	-	7,7,7	0.56	0	6,6,6	0.15	0
4	SWA	B	1909	3	13,13,13	0.81	0	13,19,19	1.60	1 (7%)

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	PG0	A	1902	-	-	0/5/5/5	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PG0	A	1903	-	-	0/5/5/5	0/0/0/0
2	PG0	A	1904	-	-	0/5/5/5	0/0/0/0
2	PG0	A	1905	-	-	0/5/5/5	0/0/0/0
4	SWA	A	1907	3	-	0/0/26/26	0/2/2/2
2	PG0	B	1901	-	-	0/5/5/5	0/0/0/0
2	PG0	B	1902	-	-	0/5/5/5	0/0/0/0
2	PG0	B	1903	-	-	0/5/5/5	0/0/0/0
2	PG0	B	1904	-	-	0/5/5/5	0/0/0/0
2	PG0	B	1905	-	-	0/5/5/5	0/0/0/0
2	PG0	B	1906	-	-	0/5/5/5	0/0/0/0
2	PG0	B	1907	-	-	0/5/5/5	0/0/0/0
4	SWA	B	1909	3	-	0/0/26/26	1/2/2/2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1907	SWA	C5-N4-C3	-3.75	109.04	111.85
4	B	1909	SWA	C5-N4-C3	4.89	115.50	111.85

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1909	SWA	C1-C2-C3-C5-C6-N4

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1905	PG0	1	0
4	A	1907	SWA	1	0
2	B	1902	PG0	2	0
2	B	1903	PG0	2	0
2	B	1905	PG0	1	0
4	B	1909	SWA	3	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 ⓘ

### 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	896/923 (97%)	-0.24	10 (1%)	82 79	17, 26, 41, 66	0
1	B	896/923 (97%)	-0.20	11 (1%)	81 77	18, 27, 41, 66	0
All	All	1792/1846 (97%)	-0.22	21 (1%)	81 77	17, 27, 41, 66	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	162	THR	4.0
1	A	162	THR	3.8
1	B	521	THR	3.6
1	B	520	GLU	3.6
1	B	3	THR	3.4
1	A	159	GLN	3.0
1	A	156	PHE	2.7
1	A	160	VAL	2.7
1	A	157	ASN	2.6
1	A	256	PHE	2.5
1	A	207	LEU	2.5
1	B	901	GLN	2.3
1	B	156	PHE	2.3
1	A	249	GLN	2.3
1	B	522	VAL	2.2
1	B	225	LEU	2.2
1	B	271	ASP	2.1
1	A	2	ALA	2.1
1	B	269	ALA	2.0
1	A	158	ASN	2.0
1	B	45	ASP	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no 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	PG0	A	1904	8/8	0.93	0.27	9.82	42,45,47,47	0
2	PG0	B	1903	8/8	0.87	0.49	8.62	66,67,67,68	0
2	PG0	B	1904	8/8	0.83	0.33	7.13	79,81,85,85	0
4	SWA	B	1909	12/12	0.91	0.28	7.09	45,49,50,51	0
2	PG0	A	1905	8/8	0.89	0.34	6.47	44,50,54,55	0
2	PG0	A	1902	8/8	0.95	0.28	4.93	47,48,50,51	0
2	PG0	B	1902	8/8	0.93	0.25	4.85	45,48,52,53	0
4	SWA	A	1907	12/12	0.93	0.23	4.25	44,49,50,51	0
2	PG0	B	1906	8/8	0.68	0.33	3.86	77,81,84,84	0
3	ZN	B	1908	1/1	0.98	0.15	0.80	70,70,70,70	0
2	PG0	B	1901	8/8	0.74	0.19	0.26	82,83,83,84	0
3	ZN	A	1906	1/1	0.92	0.09	-2.06	69,69,69,69	0
2	PG0	A	1903	8/8	0.95	0.23	-	51,52,53,54	0
2	PG0	B	1907	8/8	0.72	0.28	-	91,95,97,97	0
2	PG0	B	1905	8/8	0.83	0.35	-	92,93,93,93	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.