Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 06:18 AM GMT

PDB ID	:	2WMJ
Title	:	CRYSTAL STRUCTURE OF THE CATALYTIC MODULE OF A FAMILY 98 GLYCOSIDE HYDROLASE FROM STREPTOCOCCUS PNEUMONIAE SP3-BS71 (SP3GH98) IN COMPLEX WITH THE B-TRISACCHARIDE BLOOD GROUP ANTIGEN.
Authors	:	Higgins, M.A.; Whitworth, G.E.; El Warry, N.; Randriantsoa, M.; Samain, E.; Burke, R.D.; Vocadlo, D.J.; Boraston, A.B.
Deposited on	:	2009-06-30
Resolution	:	2.00 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

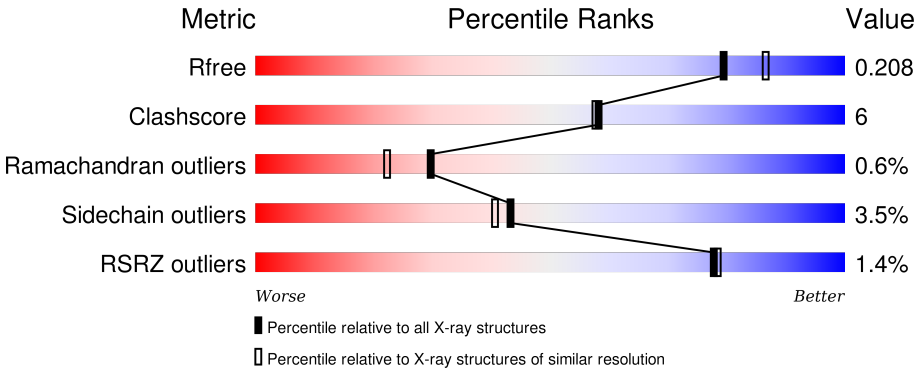
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 >=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 <=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	606	% 80% 13% • 6%
1	B	606	% 82% 11% • 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10400 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 FUCOLECTIN-RELATED PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	572	Total	C	N	O	S	0	0	0
			4681	3009	783	876	13			
1	B	571	Total	C	N	O	S	0	0	0
			4673	3005	782	873	13			

- Molecule 2 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	3	Total	C	O	0	0
			33	18	15		
2	B	3	Total	C	O	0	0
			33	18	15		

- Molecule 3 is water.

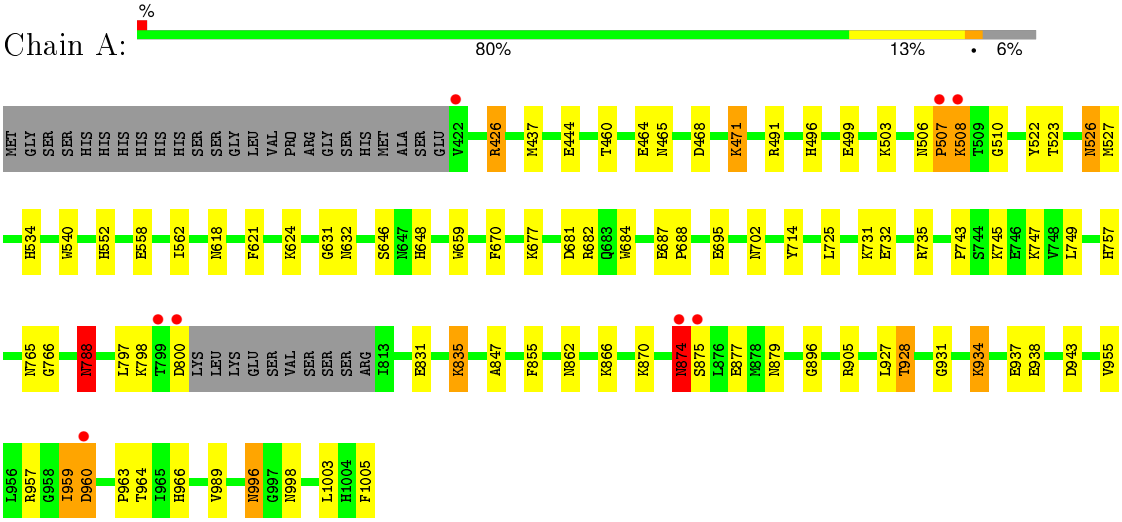
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	450	Total	O	0	0
			450	450		
3	B	530	Total	O	0	0
			530	530		

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.

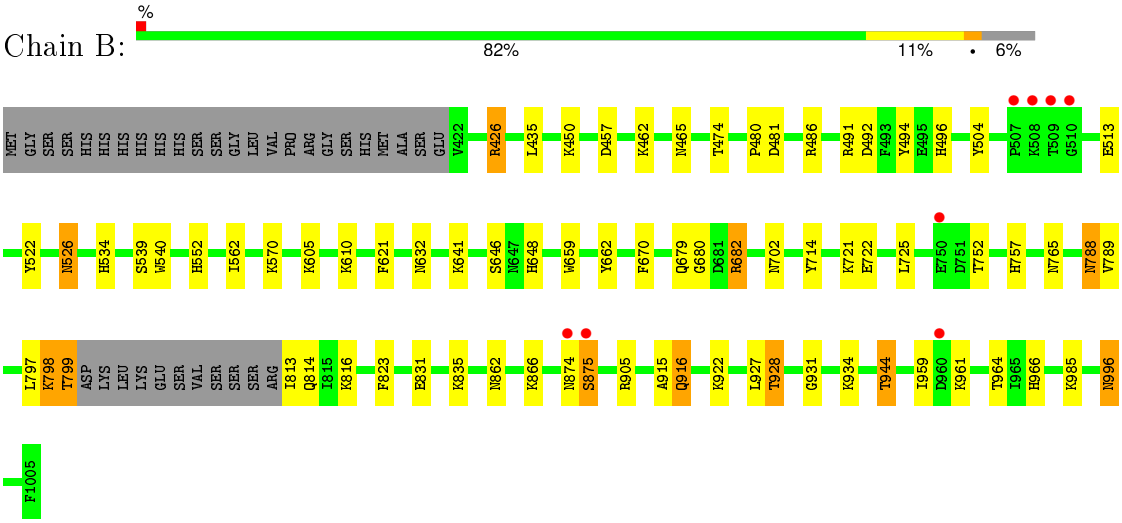
- Molecule 1: FUCOLECTIN-RELATED PROTEIN

Chain A:



- Molecule 1: FUCOLECTIN-RELATED PROTEIN

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	98.07Å 154.18Å 91.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.94 – 2.00 39.92 – 2.00	Depositor EDS
% Data completeness (in resolution range)	84.0 (39.94-2.00) 84.0 (39.92-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.74 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.161 , 0.211 0.161 , 0.208	Depositor DCC
R_{free} test set	3942 reflections (5.26%)	DCC
Wilson B-factor (Å^2)	13.8	Xtriage
Anisotropy	0.243	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 47.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	2 of 78859 reflections (0.003%)	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	10400	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.32 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.3342e-03.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GLA, GAL, FUC

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	0.74	0/4813	0.70	6/6520 (0.1%)
1	B	0.76	0/4805	0.74	7/6509 (0.1%)
All	All	0.75	0/9618	0.72	13/13029 (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	A	0	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	426	ARG	NE-CZ-NH2	-10.82	114.89	120.30
1	B	682	ARG	NE-CZ-NH1	-8.85	115.87	120.30
1	B	875	SER	N-CA-C	-7.89	89.69	111.00
1	A	426	ARG	NE-CZ-NH2	-7.33	116.63	120.30
1	B	426	ARG	NE-CZ-NH1	7.32	123.96	120.30
1	A	682	ARG	NE-CZ-NH1	-6.45	117.08	120.30
1	A	508	LYS	N-CA-C	6.13	127.56	111.00
1	A	426	ARG	NE-CZ-NH1	6.01	123.31	120.30
1	B	492	ASP	CB-CG-OD1	5.97	123.67	118.30
1	B	486	ARG	NE-CZ-NH2	-5.93	117.34	120.30
1	B	457	ASP	CB-CG-OD1	5.73	123.46	118.30
1	A	471	LYS	CD-CE-NZ	5.67	124.74	111.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	874	ASN	CB-CA-C	5.12	120.63	110.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	507	PRO	Peptide
1	A	874	ASN	Peptide
1	B	798	LYS	Peptide
1	B	874	ASN	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	4681	0	4522	61	0
1	B	4673	0	4518	48	0
2	A	33	0	30	1	0
2	B	33	0	30	1	0
3	A	450	0	0	12	2
3	B	530	0	0	12	2
All	All	10400	0	9100	109	2

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 6.

All (109) 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:721:LYS:HE3	3:B:3270:HOH:O	1.53	1.09
1:B:799:THR:HB	3:B:3352:HOH:O	1.64	0.97
1:B:927:LEU:O	1:B:928:THR:HB	1.70	0.89
1:A:928:THR:HG22	1:A:931:GLY:H	1.37	0.87
1:B:928:THR:HG22	1:B:931:GLY:H	1.36	0.86
1:B:831:GLU:O	1:B:835:LYS:HE2	1.81	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:757:HIS:HE1	1:A:797:LEU:O	1.65	0.80
1:A:905:ARG:H	1:A:996:ASN:HD21	1.29	0.79
1:A:670:PHE:H	1:A:862:ASN:HD21	1.31	0.78
1:A:934:LYS:HE3	1:A:937:GLU:OE1	1.86	0.76
1:B:905:ARG:H	1:B:996:ASN:HD21	1.30	0.75
1:B:496:HIS:HD2	3:B:3024:HOH:O	1.70	0.75
1:A:465:ASN:HB3	3:A:3226:HOH:O	1.88	0.73
1:B:757:HIS:HE1	1:B:797:LEU:O	1.72	0.71
1:A:558:GLU:HB2	1:A:624:LYS:HE3	1.73	0.71
1:B:670:PHE:H	1:B:862:ASN:HD21	1.39	0.70
1:B:462:LYS:HE3	3:B:3038:HOH:O	1.92	0.68
1:B:670:PHE:H	1:B:862:ASN:ND2	1.92	0.67
1:A:632:ASN:ND2	1:A:765:ASN:HB3	2.11	0.65
1:A:927:LEU:O	1:A:928:THR:HB	1.95	0.65
1:A:426:ARG:HE	1:A:552:HIS:HD2	1.42	0.64
1:B:757:HIS:CE1	1:B:797:LEU:O	2.51	0.63
1:A:670:PHE:H	1:A:862:ASN:ND2	1.95	0.63
1:A:964:THR:OG1	1:A:966:HIS:HE1	1.82	0.63
1:A:959:ILE:HG21	1:A:989:VAL:HG23	1.82	0.62
1:B:702:ASN:HD21	1:B:788:ASN:HB2	1.66	0.60
1:A:1005:PHE:HA	3:A:3443:HOH:O	2.00	0.60
1:A:732:GLU:HG2	3:A:3233:HOH:O	2.01	0.60
1:A:444:GLU:HG3	3:A:3044:HOH:O	2.03	0.59
1:B:799:THR:HG22	3:B:3353:HOH:O	2.01	0.59
1:B:964:THR:OG1	1:B:966:HIS:HE1	1.86	0.58
1:B:916:GLN:N	1:B:916:GLN:HE21	2.01	0.58
1:B:944:THR:CG2	3:B:3299:HOH:O	2.51	0.57
1:A:526:ASN:O	1:A:534:HIS:HD2	1.88	0.57
1:A:702:ASN:HD21	1:A:788:ASN:HB2	1.71	0.56
1:B:646:SER:OG	1:B:648:HIS:HD2	1.88	0.56
1:A:646:SER:OG	1:A:648:HIS:HD2	1.88	0.55
1:A:499:GLU:O	1:A:503:LYS:HG2	2.07	0.55
1:A:659:TRP:CE3	2:A:2007:GAL:H62	2.42	0.55
1:B:659:TRP:CE3	2:B:2007:GAL:H62	2.42	0.54
1:A:870:LYS:NZ	1:A:877:GLU:OE2	2.34	0.54
1:A:934:LYS:HE2	1:A:938:GLU:HG3	1.88	0.54
1:A:757:HIS:CE1	1:A:797:LEU:O	2.54	0.53
1:A:747:LYS:HE3	3:A:3316:HOH:O	2.09	0.53
1:B:526:ASN:O	1:B:534:HIS:HD2	1.91	0.53
1:B:522:TYR:HA	1:B:526:ASN:ND2	2.25	0.52
1:A:468:ASP:OD1	1:A:471:LYS:HE2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:788:ASN:HD22	1:B:789:VAL:H	1.58	0.51
1:A:731:LYS:HD3	1:A:735:ARG:NH1	2.25	0.51
1:B:450:LYS:HD3	3:B:3019:HOH:O	2.10	0.51
1:B:522:TYR:HA	1:B:526:ASN:HD21	1.76	0.51
1:A:491:ARG:HD2	1:A:540:TRP:CD1	2.46	0.50
1:A:496:HIS:HD2	3:A:3018:HOH:O	1.95	0.50
1:B:816:LYS:CB	1:B:823:PHE:HE2	2.25	0.50
1:A:964:THR:OG1	1:A:966:HIS:CE1	2.63	0.50
1:B:928:THR:HG23	3:B:3453:HOH:O	2.11	0.49
1:A:1005:PHE:HB2	3:A:3403:HOH:O	2.12	0.49
1:A:963:PRO:HD3	1:A:989:VAL:HG21	1.95	0.49
1:B:632:ASN:ND2	1:B:765:ASN:HB3	2.25	0.49
1:B:570:LYS:HE2	3:B:3128:HOH:O	2.12	0.49
1:B:641:LYS:NZ	1:B:702:ASN:HD22	2.11	0.49
1:A:632:ASN:HD21	1:A:765:ASN:HB3	1.76	0.49
1:B:944:THR:HG23	3:B:3299:HOH:O	2.12	0.48
1:A:527:MET:HE2	3:A:3075:HOH:O	2.13	0.48
1:A:631:GLY:O	1:A:766:GLY:HA3	2.14	0.48
1:B:632:ASN:HD21	1:B:765:ASN:HB3	1.79	0.48
1:A:896:GLY:HA3	1:A:1003:LEU:O	2.12	0.47
1:A:522:TYR:HA	1:A:526:ASN:ND2	2.30	0.47
1:A:963:PRO:CD	1:A:989:VAL:HG21	2.44	0.47
1:A:526:ASN:O	1:A:534:HIS:CD2	2.68	0.47
1:B:788:ASN:HD22	1:B:789:VAL:N	2.14	0.46
1:B:922:LYS:HE2	3:B:3231:HOH:O	2.15	0.46
1:A:960:ASP:OD1	1:A:960:ASP:N	2.37	0.46
1:B:504:TYR:HD2	1:B:513:GLU:O	1.99	0.46
1:A:522:TYR:HA	1:A:526:ASN:HD21	1.80	0.46
1:A:905:ARG:H	1:A:996:ASN:ND2	2.07	0.45
1:A:465:ASN:CG	3:A:3027:HOH:O	2.54	0.45
1:A:503:LYS:HB3	1:A:503:LYS:HE2	1.81	0.45
1:A:426:ARG:HE	1:A:552:HIS:CD2	2.29	0.45
1:A:681:ASP:HA	1:A:684:TRP:NE1	2.32	0.45
1:B:435:LEU:HB3	1:B:474:THR:HG22	1.98	0.44
1:B:610:LYS:HD2	3:B:3168:HOH:O	2.17	0.44
1:B:816:LYS:HB3	1:B:823:PHE:HE2	1.81	0.44
1:A:618:ASN:ND2	3:A:3151:HOH:O	2.51	0.44
1:A:714:TYR:CZ	1:A:725:LEU:HD12	2.53	0.44
1:B:426:ARG:HH21	1:B:552:HIS:HD2	1.66	0.43
1:A:460:THR:O	1:A:464:GLU:HG3	2.17	0.43
1:A:847:ALA:HA	1:A:855:PHE:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:959:ILE:HG21	1:A:989:VAL:CG2	2.47	0.43
1:B:662:TYR:CD1	1:B:680:GLY:HA2	2.53	0.43
1:B:480:PRO:HD3	1:B:494:TYR:OH	2.18	0.43
1:A:659:TRP:CZ2	1:A:684:TRP:CD1	3.06	0.43
1:B:465:ASN:HD21	1:B:722:GLU:HG3	1.84	0.43
1:B:714:TYR:CZ	1:B:725:LEU:HD12	2.54	0.43
1:B:964:THR:OG1	1:B:966:HIS:CE1	2.70	0.43
1:B:641:LYS:NZ	1:B:788:ASN:ND2	2.67	0.43
1:A:506:ASN:O	1:A:510:GLY:N	2.52	0.43
1:A:831:GLU:O	1:A:835:LYS:HG2	2.20	0.42
1:A:866:LYS:HE3	3:A:3346:HOH:O	2.19	0.42
1:B:915:ALA:C	1:B:916:GLN:HE21	2.23	0.42
1:B:679:GLN:OE1	1:B:682:ARG:NH1	2.53	0.42
1:A:745:LYS:O	1:A:749:LEU:HB2	2.19	0.41
1:B:491:ARG:HD2	1:B:540:TRP:CD1	2.56	0.41
1:A:687:GLU:OE2	1:A:695:GLU:OE2	2.39	0.41
1:A:523:THR:H	1:A:526:ASN:ND2	2.18	0.40
1:B:752:THR:O	1:B:813:ILE:HG12	2.21	0.40
1:A:677:LYS:HE3	3:A:3180:HOH:O	2.21	0.40
1:A:879:ASN:HB2	1:A:955:VAL:HB	2.03	0.40
1:A:743:PRO:HB3	1:A:747:LYS:HD3	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:3271:HOH:O	3:B:3067:HOH:O[3_554]	1.55	0.65
3:A:3271:HOH:O	3:B:3068:HOH:O[3_554]	2.18	0.02

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	568/606 (94%)	540 (95%)	23 (4%)	5 (1%)	21 13
1	B	567/606 (94%)	542 (96%)	23 (4%)	2 (0%)	39 33
All	All	1135/1212 (94%)	1082 (95%)	46 (4%)	7 (1%)	30 22

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	507	PRO
1	A	508	LYS
1	A	874	ASN
1	B	481	ASP
1	A	788	ASN
1	B	562	ILE
1	A	562	ILE

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	502/532 (94%)	485 (97%)	17 (3%)	44 41
1	B	501/532 (94%)	483 (96%)	18 (4%)	42 39
All	All	1003/1064 (94%)	968 (96%)	35 (4%)	43 40

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	437	MET
1	A	526	ASN
1	A	621	PHE
1	A	688	PRO
1	A	788	ASN
1	A	798	LYS
1	A	800	ASP
1	A	835	LYS
1	A	875	SER

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Mol	Chain	Res	Type
1	A	928	THR
1	A	934	LYS
1	A	943	ASP
1	A	957	ARG
1	A	959	ILE
1	A	960	ASP
1	A	996	ASN
1	A	998	ASN
1	B	526	ASN
1	B	539	SER
1	B	605	LYS
1	B	621	PHE
1	B	788	ASN
1	B	798	LYS
1	B	799	THR
1	B	814	GLN
1	B	866	LYS
1	B	875	SER
1	B	916	GLN
1	B	928	THR
1	B	934	LYS
1	B	944	THR
1	B	959	ILE
1	B	961	LYS
1	B	985	LYS
1	B	996	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	448	GLN
1	A	465	ASN
1	A	496	HIS
1	A	512	ASN
1	A	526	ASN
1	A	534	HIS
1	A	552	HIS
1	A	581	ASN
1	A	609	GLN
1	A	618	ASN
1	A	632	ASN
1	A	648	HIS

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Mol	Chain	Res	Type
1	A	702	ASN
1	A	740	HIS
1	A	757	HIS
1	A	765	ASN
1	A	771	ASN
1	A	788	ASN
1	A	862	ASN
1	A	966	HIS
1	A	973	ASN
1	A	987	HIS
1	A	994	ASN
1	A	996	ASN
1	A	998	ASN
1	B	465	ASN
1	B	496	HIS
1	B	512	ASN
1	B	526	ASN
1	B	534	HIS
1	B	552	HIS
1	B	618	ASN
1	B	632	ASN
1	B	648	HIS
1	B	702	ASN
1	B	740	HIS
1	B	757	HIS
1	B	771	ASN
1	B	788	ASN
1	B	862	ASN
1	B	900	ASN
1	B	916	GLN
1	B	919	ASN
1	B	966	HIS
1	B	986	ASN
1	B	994	ASN
1	B	996	ASN
1	B	998	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)

6 carbohydrates 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	GLA	A	2006	2	11,11,12	0.57	0	14,15,17	0.93	0
2	GAL	A	2007	2	12,12,12	0.73	0	17,17,17	1.44	4 (23%)
2	FUC	A	2008	2	10,10,11	0.74	0	14,14,16	1.20	1 (7%)
2	GLA	B	2006	2	11,11,12	0.71	0	14,15,17	0.85	0
2	GAL	B	2007	2	12,12,12	0.74	0	17,17,17	1.32	3 (17%)
2	FUC	B	2008	2	10,10,11	0.77	0	14,14,16	1.16	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	GLA	A	2006	2	-	0/2/19/22	0/1/1/1
2	GAL	A	2007	2	-	0/2/22/22	0/1/1/1
2	FUC	A	2008	2	-	0/0/17/20	0/1/1/1
2	GLA	B	2006	2	-	0/2/19/22	0/1/1/1
2	GAL	B	2007	2	-	0/2/22/22	0/1/1/1
2	FUC	B	2008	2	-	0/0/17/20	0/1/1/1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	2007	GAL	O3-C3-C2	-3.13	103.30	110.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	2007	GAL	O1-C1-O5	-2.77	102.68	110.25
2	B	2007	GAL	O1-C1-O5	-2.43	103.59	110.25
2	B	2007	GAL	O3-C3-C2	-2.16	105.47	110.34
2	A	2007	GAL	O5-C1-C2	2.02	113.03	109.80
2	B	2008	FUC	O5-C5-C6	2.11	109.62	106.13
2	B	2007	GAL	O5-C1-C2	2.23	113.36	109.80
2	A	2007	GAL	C1-C2-C3	2.39	113.98	110.43
2	A	2008	FUC	C1-O5-C5	3.20	117.32	112.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2007	GAL	1	0
2	B	2007	GAL	1	0

5.6 Ligand geometry (i)

There are no ligands in this entry.

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^{th} 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	572/606 (94%)	-0.39	8 (1%) 78 78	4, 11, 29, 42	0
1	B	571/606 (94%)	-0.43	8 (1%) 78 78	3, 10, 24, 43	0
All	All	1143/1212 (94%)	-0.41	16 (1%) 78 78	3, 10, 26, 43	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	507	PRO	4.2
1	A	508	LYS	4.0
1	A	507	PRO	3.4
1	A	875	SER	3.3
1	B	509	THR	3.0
1	A	422	VAL	2.9
1	B	510	GLY	2.7
1	B	960	ASP	2.7
1	A	960	ASP	2.6
1	B	508	LYS	2.6
1	A	800	ASP	2.5
1	A	874	ASN	2.2
1	B	750	GLU	2.1
1	B	874	ASN	2.1
1	A	799	THR	2.0
1	B	875	SER	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 (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^{th} 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^2)	Q<0.9
2	FUC	A	2008	10/11	0.99	0.10	0.55	5,8,9,9	0
2	FUC	B	2008	10/11	0.98	0.11	0.48	6,7,8,9	0
2	GAL	B	2007	12/12	0.97	0.09	-0.03	6,9,12,17	0
2	GAL	A	2007	12/12	0.95	0.08	-0.34	6,10,13,17	0
2	GLA	B	2006	11/12	0.99	0.10	-0.47	3,4,6,6	0
2	GLA	A	2006	11/12	0.99	0.08	-1.53	5,6,8,9	0

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.