



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2017 – 01:29 PM EST

PDB ID : 2WGC
Title : 2.2 ANGSTROMS RESOLUTION STRUCTURE ANALYSIS OF TWO RE-
FINED N-ACETYLNEURAMINYLLACTOSE-WHEAT GERM AGGLU-
TININ ISOLECTIN COMPLEXES
Authors : Wright, C.S.
Deposited on : unknown
Resolution : 2.20 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

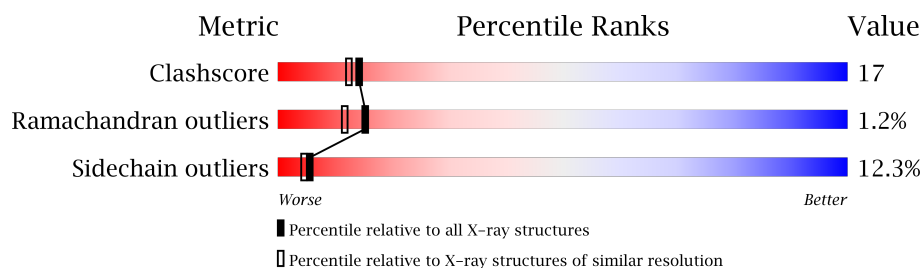
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 $\leq 5\%$. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	171	 63% 29% 6% •
1	B	171	 61% 32% 6% •

2 Entry composition [i](#)

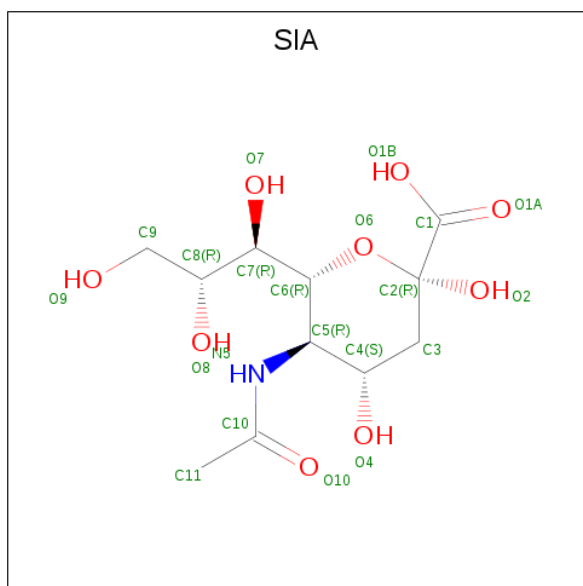
There are 5 unique types of molecules in this entry. The entry contains 2634 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 WHEAT GERM LECTIN.

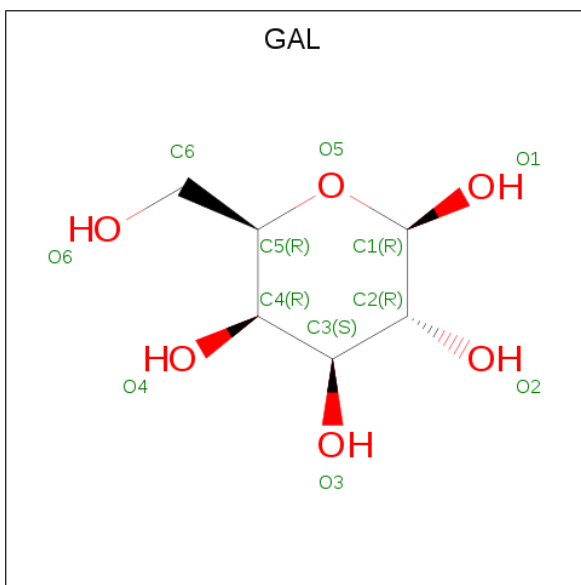
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	0	0	0
			1159	674	215	236	34			
1	B	171	Total	C	N	O	S	0	0	0
			1154	669	215	236	34			

- Molecule 2 is O-SIALIC ACID (three-letter code: SIA) (formula: $C_{11}H_{19}NO_9$).



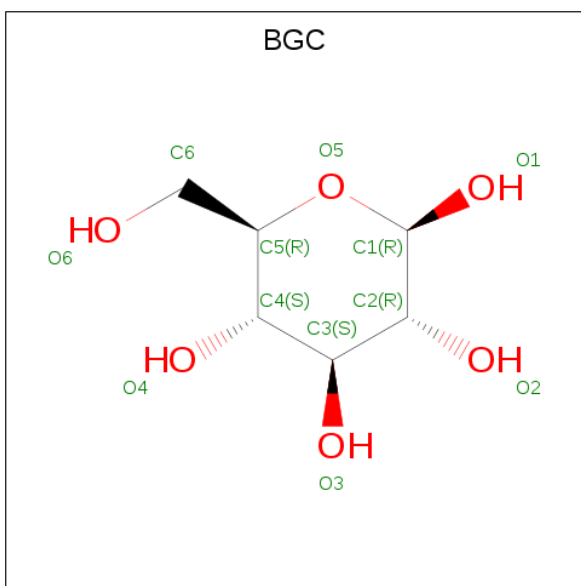
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			21	11	1	9		
2	B	1	Total	C	N	O	0	0
			21	11	1	9		

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



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

- Molecule 4 is BETA-D-GLUCOSE (three-letter code: BGC) (formula: $C_6H_{12}O_6$).



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

- Molecule 5 is water.

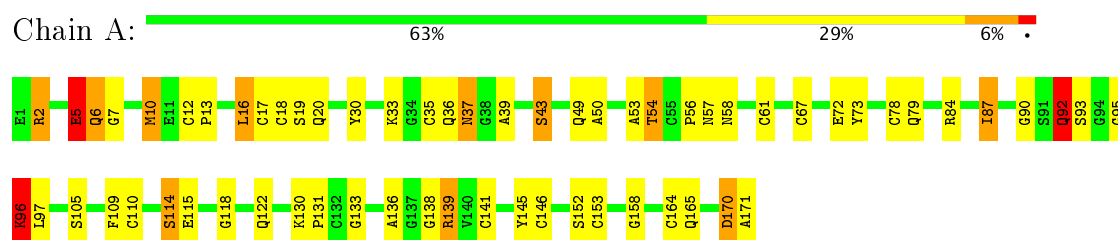
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	115	Total 115	O 115	0	0
5	B	120	Total 120	O 120	0	0

3 Residue-property plots [i](#)

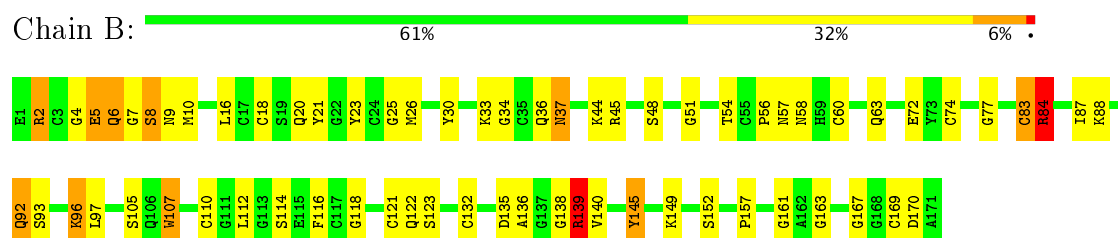
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

• Molecule 1: WHEAT GERM LECTIN



• Molecule 1: WHEAT GERM LECTIN



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	51.44Å 73.35Å 91.68Å 90.00° 97.52° 90.00°	Depositor
Resolution (Å)	8.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.153 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2634	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SIA, BGC, PCA, GAL

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	1.12	4/1175 (0.3%)	1.73	16/1575 (1.0%)
1	B	1.13	4/1168 (0.3%)	1.81	22/1563 (1.4%)
All	All	1.13	8/2343 (0.3%)	1.77	38/3138 (1.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	5	GLU	CD-OE2	7.35	1.33	1.25
1	A	72	GLU	CD-OE2	7.32	1.33	1.25
1	B	72	GLU	CD-OE2	6.79	1.33	1.25
1	B	5	GLU	CD-OE2	6.70	1.33	1.25
1	B	83	CYS	CB-SG	-5.76	1.72	1.81
1	A	115	GLU	CD-OE2	5.56	1.31	1.25
1	B	169	CYS	CB-SG	-5.32	1.73	1.81
1	A	13	PRO	N-CD	5.22	1.55	1.47

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	84	ARG	CD-NE-CZ	15.61	145.45	123.60
1	B	84	ARG	NE-CZ-NH1	14.87	127.73	120.30
1	B	139	ARG	NE-CZ-NH1	-9.15	115.73	120.30
1	B	21	TYR	CB-CG-CD2	-8.66	115.80	121.00
1	B	45	ARG	NE-CZ-NH1	8.41	124.50	120.30
1	A	78	CYS	CA-CB-SG	8.17	128.70	114.00
1	A	96	LYS	CA-CB-CG	8.09	131.19	113.40
1	A	2	ARG	NE-CZ-NH2	-7.38	116.61	120.30
1	B	45	ARG	CD-NE-CZ	7.36	133.90	123.60
1	A	96	LYS	N-CA-CB	7.24	123.62	110.60
1	B	139	ARG	NH1-CZ-NH2	7.10	127.21	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	84	ARG	NE-CZ-NH2	7.00	123.80	120.30
1	A	170	ASP	CB-CG-OD1	6.85	124.47	118.30
1	A	170	ASP	CB-CG-OD2	-6.85	112.14	118.30
1	B	2	ARG	NE-CZ-NH1	6.80	123.70	120.30
1	A	139	ARG	NE-CZ-NH2	-6.58	117.01	120.30
1	B	21	TYR	CB-CG-CD1	6.53	124.92	121.00
1	B	139	ARG	NE-CZ-NH2	-6.47	117.06	120.30
1	B	139	ARG	CA-CB-CG	6.39	127.46	113.40
1	A	170	ASP	N-CA-CB	6.29	121.93	110.60
1	A	73	TYR	CB-CG-CD2	-6.26	117.24	121.00
1	B	152	SER	N-CA-CB	6.20	119.81	110.50
1	B	145	TYR	CB-CG-CD2	-6.14	117.32	121.00
1	B	84	ARG	NH1-CZ-NH2	-6.10	112.69	119.40
1	A	115	GLU	CG-CD-OE1	5.88	130.07	118.30
1	A	43	SER	CB-CA-C	5.81	121.14	110.10
1	B	30	TYR	CB-CG-CD2	-5.57	117.66	121.00
1	A	54	THR	CA-CB-CG2	5.50	120.10	112.40
1	B	54	THR	CA-CB-CG2	5.29	119.81	112.40
1	B	145	TYR	CB-CG-CD1	5.24	124.15	121.00
1	A	2	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	B	169	CYS	O-C-N	5.18	130.99	122.70
1	B	60	CYS	O-C-N	5.11	130.88	122.70
1	A	90	GLY	O-C-N	5.10	130.85	122.70
1	B	170	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	B	30	TYR	O-C-N	5.07	130.81	122.70
1	B	107	TRP	CA-CB-CG	-5.04	104.13	113.70
1	A	10	MET	CA-CB-CG	5.03	121.85	113.30

There are no chirality outliers.

There are no planarity outliers.

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	1159	0	972	36	0
1	B	1154	0	975	41	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	21	0	17	0	0
2	B	21	0	17	0	0
3	A	11	0	9	0	0
3	B	11	0	9	0	0
4	A	11	0	10	0	0
4	B	11	0	10	0	0
5	A	115	0	0	5	0
5	B	120	0	0	8	0
All	All	2634	0	2019	77	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 17.

All (77) 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:92:GLN:H	1:B:92:GLN:HE21	1.25	0.82
1:B:92:GLN:HE22	1:B:122:GLN:HE22	1.31	0.77
1:B:92:GLN:H	1:B:92:GLN:NE2	1.86	0.72
1:A:6:GLN:H	1:A:6:GLN:HE21	1.37	0.71
1:B:2:ARG:O	1:B:7:GLY:HA3	1.93	0.68
1:B:37:ASN:OD1	5:B:356:HOH:O	2.14	0.66
1:A:164:CYS:HB3	1:A:171:ALA:O	1.96	0.66
1:A:6:GLN:HE22	1:A:36:GLN:HE22	1.45	0.65
1:A:54:THR:HG22	5:A:400:HOH:O	1.98	0.63
1:B:16:LEU:HD22	1:B:26:MET:HG2	1.80	0.63
1:B:20:GLN:NE2	5:B:411:HOH:O	2.33	0.62
1:B:107:TRP:CE3	1:B:107:TRP:HA	2.35	0.62
1:A:37:ASN:C	1:A:37:ASN:HD22	2.03	0.61
1:B:83:CYS:HB2	1:B:87:ILE:HD11	1.82	0.61
1:A:56:PRO:HD2	5:A:319:HOH:O	2.02	0.60
1:B:6:GLN:HE22	1:B:36:GLN:HE22	1.49	0.60
1:B:16:LEU:HD13	1:B:25:GLY:HA2	1.83	0.60
1:B:139:ARG:NH1	5:B:410:HOH:O	2.33	0.60
1:B:96:LYS:HD3	5:B:338:HOH:O	2.04	0.57
1:A:7:GLY:O	1:A:10:MET:HB2	2.04	0.56
1:A:61:CYS:O	1:A:79:GLN:N	2.33	0.56
1:B:37:ASN:ND2	5:B:343:HOH:O	2.39	0.56
1:A:2:ARG:O	1:A:7:GLY:HA3	2.06	0.55
1:B:132:CYS:HA	1:B:139:ARG:HG3	1.87	0.55
1:B:107:TRP:CA	1:B:107:TRP:CE3	2.91	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:ALA:O	1:A:53:ALA:HB2	2.09	0.53
1:A:133:GLY:O	1:A:136:ALA:N	2.34	0.53
1:B:20:GLN:HG2	5:B:392:HOH:O	2.09	0.53
1:B:135:ASP:HB2	5:B:344:HOH:O	2.08	0.52
1:A:6:GLN:NE2	1:A:6:GLN:H	2.07	0.52
1:B:20:GLN:HB3	1:B:34:GLY:HA3	1.92	0.52
1:A:158:GLY:HA3	5:A:369:HOH:O	2.10	0.51
1:A:17:CYS:HB3	1:A:35:CYS:SG	2.50	0.51
1:B:63:GLN:NE2	5:B:317:HOH:O	2.32	0.50
1:A:92:GLN:HE22	1:A:122:GLN:HE22	1.58	0.50
1:A:93:SER:O	1:A:96:LYS:HG2	2.11	0.50
1:A:92:GLN:NE2	1:A:92:GLN:H	2.10	0.50
1:A:18:CYS:HB2	1:A:37:ASN:HD21	1.78	0.49
1:B:87:ILE:HB	1:B:110:CYS:HB2	1.95	0.49
1:B:107:TRP:CD2	1:B:107:TRP:N	2.79	0.48
1:B:97:LEU:HD23	1:B:123:SER:OG	2.13	0.48
1:A:20:GLN:NE2	5:A:411:HOH:O	2.32	0.48
1:B:74:CYS:O	1:B:84:ARG:HB2	2.13	0.48
1:B:57:ASN:O	1:B:58:ASN:HB2	2.13	0.48
1:A:87:ILE:HB	1:A:110:CYS:HB2	1.96	0.47
1:A:114:SER:O	1:A:118:GLY:HA3	2.14	0.47
1:B:157:PRO:O	1:B:161:GLY:HA3	2.14	0.47
1:B:107:TRP:HE3	1:B:107:TRP:HA	1.75	0.47
1:B:63:GLN:HB3	1:B:77:GLY:HA3	1.96	0.46
1:A:43:SER:HB3	1:A:67:CYS:O	2.16	0.45
1:B:18:CYS:HB3	1:B:36:GLN:HB2	1.98	0.45
1:A:49:GLN:OE1	1:A:49:GLN:N	2.48	0.45
1:A:136:ALA:C	1:A:138:GLY:N	2.68	0.44
1:B:118:GLY:O	1:B:121:CYS:HB2	2.17	0.44
1:B:136:ALA:C	1:B:138:GLY:H	2.21	0.44
1:A:146:CYS:O	1:A:153:CYS:HA	2.18	0.44
1:A:12:CYS:HB2	1:A:16:LEU:O	2.17	0.44
1:A:57:ASN:O	1:A:58:ASN:HB2	2.16	0.44
1:A:141:CYS:HB3	1:A:145:TYR:HB2	2.00	0.43
1:B:92:GLN:NE2	1:B:92:GLN:N	2.62	0.43
1:A:130:LYS:HA	1:A:131:PRO:HD3	1.80	0.43
1:B:149:LYS:HB3	1:B:163:GLY:HA3	2.01	0.42
1:A:105:SER:HA	5:A:380:HOH:O	2.20	0.42
1:A:33:LYS:O	1:A:33:LYS:HD3	2.19	0.42
1:B:16:LEU:HD13	1:B:25:GLY:CA	2.49	0.42
1:A:92:GLN:HE21	1:A:92:GLN:H	1.66	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:SER:HB3	1:B:116:PHE:HA	2.02	0.42
1:A:19:SER:HB3	1:A:30:TYR:HA	2.03	0.41
1:B:4:GLY:O	1:B:9:ASN:N	2.50	0.41
1:B:92:GLN:N	1:B:92:GLN:HE21	2.03	0.41
1:B:2:ARG:NH1	1:B:23:TYR:OH	2.54	0.41
1:A:87:ILE:O	1:A:109:PHE:HA	2.21	0.41
1:A:95:GLY:O	1:A:96:LYS:C	2.59	0.41
1:A:5:GLU:HG3	1:A:5:GLU:O	2.20	0.41
1:B:145:TYR:O	1:B:167:GLY:HA3	2.20	0.41
1:B:56:PRO:O	1:B:57:ASN:HB2	2.20	0.41
1:B:7:GLY:C	1:B:8:SER:OG	2.60	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	169/171 (99%)	149 (88%)	18 (11%)	2 (1%)	15	12
1	B	169/171 (99%)	156 (92%)	11 (6%)	2 (1%)	15	12
All	All	338/342 (99%)	305 (90%)	29 (9%)	4 (1%)	15	12

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	93	SER
1	A	39	ALA
1	A	92	GLN
1	B	51	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	118/120 (98%)	105 (89%)	13 (11%)	7	6
1	B	118/120 (98%)	102 (86%)	16 (14%)	4	3
All	All	236/240 (98%)	207 (88%)	29 (12%)	5	4

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLU
1	A	6	GLN
1	A	16	LEU
1	A	37	ASN
1	A	87	ILE
1	A	92	GLN
1	A	96	LYS
1	A	97	LEU
1	A	114	SER
1	A	139	ARG
1	A	152	SER
1	A	165	GLN
1	A	170	ASP
1	B	5	GLU
1	B	6	GLN
1	B	8	SER
1	B	10	MET
1	B	33	LYS
1	B	37	ASN
1	B	44	LYS
1	B	48	SER
1	B	84	ARG
1	B	88	LYS
1	B	92	GLN
1	B	96	LYS
1	B	112	LEU
1	B	114	SER

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Mol	Chain	Res	Type
1	B	139	ARG
1	B	140	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	9	ASN
1	A	37	ASN
1	A	63	GLN
1	A	92	GLN
1	B	6	GLN
1	B	9	ASN
1	B	37	ASN
1	B	92	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	PCA	A	1	1	8,8,9	1.05	1 (12%)	9,10,12	2.10	3 (33%)
1	PCA	B	1	1	8,8,9	0.88	0	9,10,12	1.75	1 (11%)

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
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1
1	PCA	B	1	1	-	0/0/11/13	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	PCA	O-C	2.53	1.30	1.19

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1	PCA	O-C-CA	-4.79	113.99	125.15
1	A	1	PCA	O-C-CA	-4.65	114.30	125.15
1	A	1	PCA	CB-CG-CD	-2.14	100.67	104.30
1	A	1	PCA	CB-CA-C	2.61	116.30	112.70

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 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	A	301	3	18,21,21	1.12	2 (11%)	19,31,31	1.94	5 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GAL	A	302	2,4	11,11,12	0.77	0	10,15,17	1.07	1 (10%)
4	BGC	A	303	3	11,11,12	1.29	1 (9%)	15,15,17	1.68	4 (26%)
2	SIA	B	301	3	18,21,21	0.99	0	19,31,31	1.59	4 (21%)
3	GAL	B	302	2,4	11,11,12	1.06	1 (9%)	10,15,17	1.58	1 (10%)
4	BGC	B	303	3	11,11,12	1.35	1 (9%)	15,15,17	1.33	2 (13%)

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	A	301	3	-	0/14/38/38	0/1/1/1
3	GAL	A	302	2,4	-	0/2/18/22	0/1/1/1
4	BGC	A	303	3	-	0/2/18/22	0/1/1/1
2	SIA	B	301	3	-	0/14/38/38	0/1/1/1
3	GAL	B	302	2,4	-	0/2/18/22	0/1/1/1
4	BGC	B	303	3	-	0/2/18/22	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	SIA	C3-C4	-2.68	1.49	1.53
3	B	302	GAL	C1-C2	-2.55	1.50	1.52
2	A	301	SIA	O2-C2	2.84	1.42	1.39
4	B	303	BGC	O2-C2	3.05	1.50	1.43
4	A	303	BGC	O2-C2	3.11	1.50	1.43

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	303	BGC	C3-C4-C5	-3.04	106.20	111.26
2	B	301	SIA	O10-C10-C11	-2.87	116.83	122.06
2	A	301	SIA	C11-C10-N5	-2.63	111.37	116.11
2	B	301	SIA	C6-C5-N5	-2.20	107.13	111.00
3	A	302	GAL	O1-C1-O5	-2.16	103.80	110.20
4	A	303	BGC	O5-C5-C4	-2.07	106.55	109.98
4	B	303	BGC	O3-C3-C2	-2.06	106.13	110.17
4	A	303	BGC	O5-C5-C6	2.26	110.62	106.88
2	A	301	SIA	C4-C5-N5	2.30	115.14	110.40
4	A	303	BGC	C4-C3-C2	2.56	113.94	110.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	303	BGC	C4-C3-C2	2.63	114.02	110.59
3	B	302	GAL	O5-C1-C2	2.94	114.82	110.03
2	B	301	SIA	O4-C4-C5	3.28	117.31	110.40
2	A	301	SIA	O8-C8-C9	3.40	117.06	109.21
2	B	301	SIA	O10-C10-N5	3.47	128.59	121.92
2	A	301	SIA	O10-C10-N5	4.17	129.94	121.92
2	A	301	SIA	C5-N5-C10	4.18	133.79	123.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.