



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 5, 2018 – 09:46 PM EST

PDB ID : 1OP5
Title : Crystal Structure of Fab 2G12 bound to Man9GlcNAc2
Authors : Calarese, D.A.; Scanlan, C.N.; Zwick, M.B.; Deechongkit, S.; Mimura, Y.; Kunert, R.; Stanfield, R.L.; Kelly, J.W.; Rudd, P.M.; Dwek, R.A.; Katinger, H.; Burton, D.R.; Wilson, I.A.
Deposited on : 2003-03-04
Resolution : 3.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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-20030736

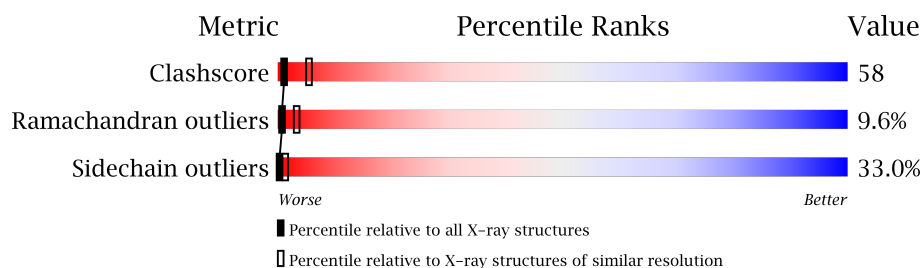
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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))
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)

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 $\leq 5\%$. 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	K	212	
1	L	212	
2	H	225	
2	M	225	

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
3	MAN	H	234	-	-	X	-
3	MAN	H	237	-	-	X	-
3	MAN	K	218	-	-	X	-
3	MAN	L	214	-	-	X	-
3	MAN	M	230	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6848 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.

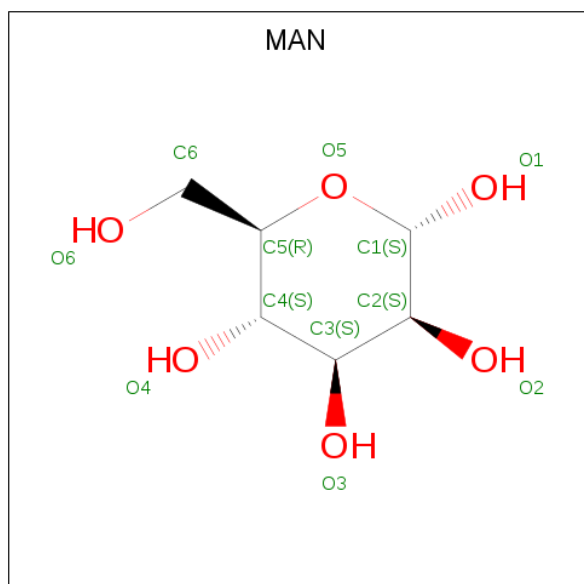
- Molecule 1 is a protein called FAB 2G12, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	212	Total	C	N	O	S	0	0	0
			1620	1016	273	326	5			
1	K	212	Total	C	N	O	S	0	0	0
			1620	1016	273	326	5			

- Molecule 2 is a protein called FAB 2G12, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	225	Total	C	N	O	S	0	0	0
			1676	1051	287	331	7			
2	M	225	Total	C	N	O	S	0	0	0
			1676	1051	287	331	7			

- Molecule 3 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula: C₆H₁₂O₆).



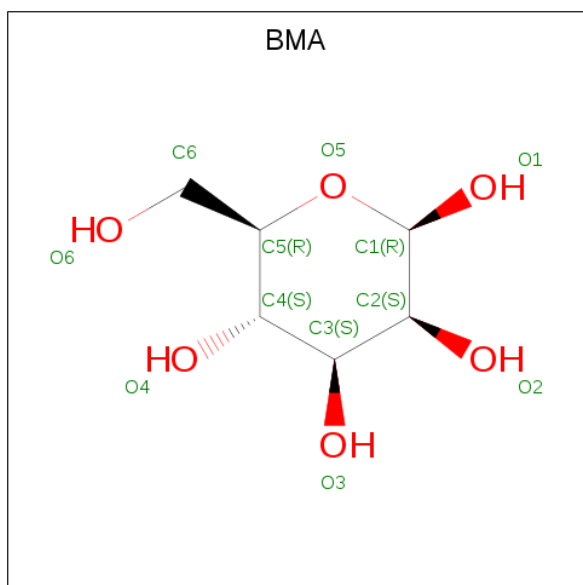
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	L	1	Total	C	O	0	0
			11	6	5		
3	L	1	Total	C	O	0	0
			11	6	5		
3	H	1	Total	C	O	0	0
			11	6	5		
3	H	1	Total	C	O	0	0
			11	6	5		
3	H	1	Total	C	O	0	0
			11	6	5		
3	H	1	Total	C	O	0	0
			11	6	5		
3	H	1	Total	C	O	0	0
			11	6	5		
3	K	1	Total	C	O	0	0
			11	6	5		
3	K	1	Total	C	O	0	0
			11	6	5		
3	K	1	Total	C	O	0	0
			11	6	5		
3	M	1	Total	C	O	0	0
			11	6	5		
3	M	1	Total	C	O	0	0
			11	6	5		
3	M	1	Total	C	O	0	0
			11	6	5		
3	M	1	Total	C	O	0	0
			11	6	5		

- Molecule 4 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	H	1	Total	C	N	O	0	0
			15	8	1	6		
4	H	1	Total	C	N	O	0	0
			14	8	1	5		
4	K	1	Total	C	N	O	0	0
			15	8	1	6		
4	K	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is BETA-D-MANNOSE (three-letter code: BMA) (formula: $C_6H_{12}O_6$).



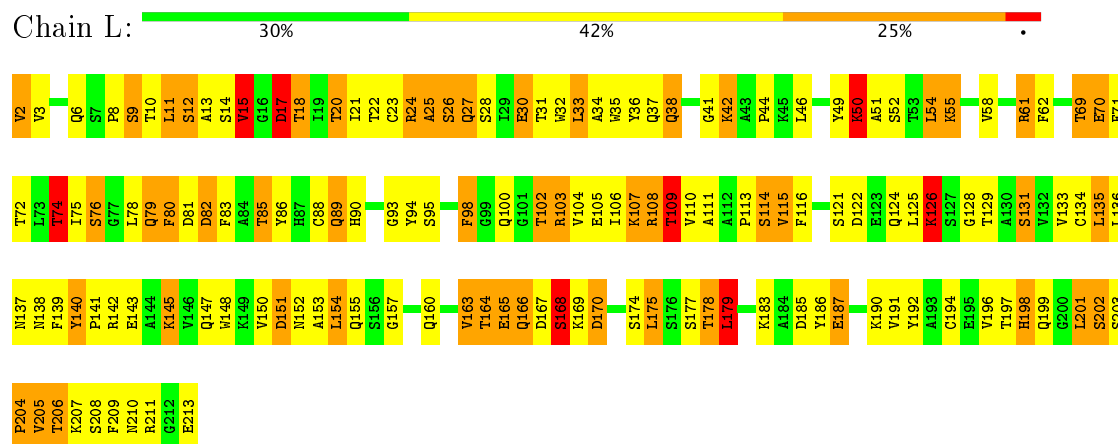
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	H	1	Total	C	O	0	0
			11	6	5		
5	K	1	Total	C	O	0	0
			11	6	5		

3 Residue-property plots

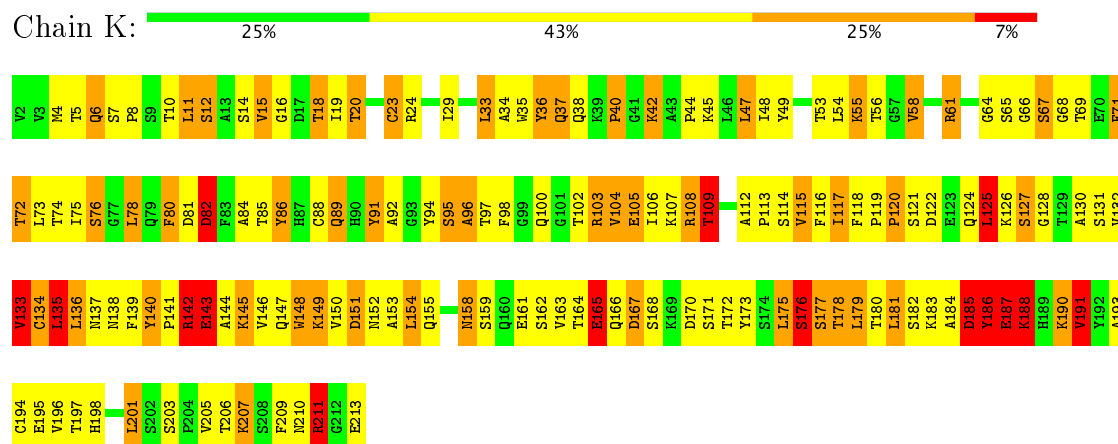
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

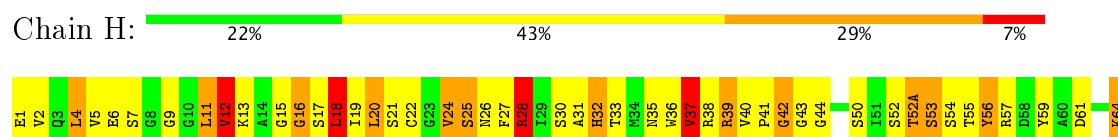
- Molecule 1: FAB 2G12, light chain

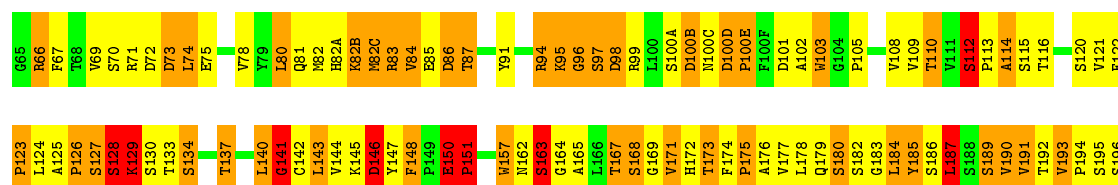


- Molecule 1: FAB 2G12, light chain

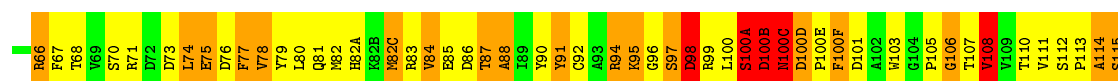
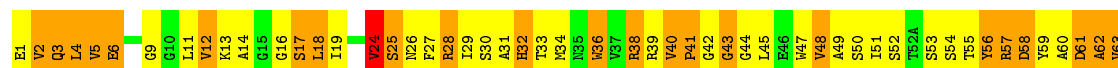
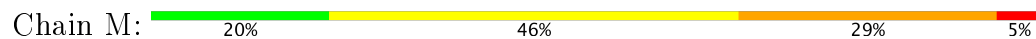


- Molecule 2: FAB 2G12, heavy chain





- Molecule 2: FAB 2G12, heavy chain



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	135.82Å 145.71Å 148.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	105.41 – 3.00	Depositor
% Data completeness (in resolution range)	97.5 (105.41-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.248 , 0.329	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6848	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, MAN

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	K	1.49	15/1656 (0.9%)	1.47	14/2251 (0.6%)
1	L	1.75	22/1656 (1.3%)	1.62	26/2251 (1.2%)
2	H	2.15	34/1715 (2.0%)	1.72	35/2337 (1.5%)
2	M	1.54	17/1714 (1.0%)	1.66	28/2333 (1.2%)
All	All	1.75	88/6741 (1.3%)	1.62	103/9172 (1.1%)

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	K	0	1
1	L	0	1
2	H	0	4
2	M	0	9
All	All	0	15

The worst 5 of 88 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	222	LYS	CE-NZ	41.93	2.53	1.49
2	H	112	SER	C-N	19.01	1.70	1.34
2	H	229	SER	CB-OG	16.80	1.64	1.42
2	H	228	LYS	CE-NZ	14.28	1.84	1.49
2	H	128	SER	CB-OG	12.46	1.58	1.42

The worst 5 of 103 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	98	ASP	CB-CG-OD2	10.81	128.03	118.30
2	M	96	GLY	N-CA-C	10.18	138.55	113.10
2	M	100(B)	ASP	CB-CG-OD2	10.05	127.34	118.30
2	M	61	ASP	CB-CG-OD2	9.67	127.01	118.30
1	K	167	ASP	CB-CG-OD2	9.45	126.80	118.30

There are no chirality outliers.

5 of 15 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	128	SER	Peptide
2	H	141	GLY	Peptide
2	H	42	GLY	Peptide
2	H	96	GLY	Peptide
1	L	25	ALA	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	K	1620	0	1551	199	0
1	L	1620	0	1551	181	0
2	H	1676	0	1628	184	0
2	M	1676	0	1628	198	0
3	H	66	0	55	20	0
3	K	33	0	29	15	0
3	L	22	0	19	6	0
3	M	55	0	48	18	0
4	H	29	0	26	0	0
4	K	29	0	26	0	0
5	H	11	0	8	4	0
5	K	11	0	8	3	0
All	All	6848	0	6577	771	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 58.

The worst 5 of 771 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:L:2:VAL:CB	1:L:2:VAL:CG2	1.77	1.58
2:M:82:MET:CE	2:M:82:MET:SD	2.01	1.48
2:H:112:SER:C	2:H:113:PRO:N	1.70	1.45
1:K:113:PRO:HD3	1:K:198:HIS:CD2	1.50	1.44
2:H:229:SER:OG	2:H:229:SER:CB	1.64	1.42

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	K	210/212 (99%)	158 (75%)	39 (19%)	13 (6%)	2	10
1	L	210/212 (99%)	177 (84%)	24 (11%)	9 (4%)	3	18
2	H	223/225 (99%)	158 (71%)	42 (19%)	23 (10%)	0	3
2	M	221/225 (98%)	155 (70%)	28 (13%)	38 (17%)	0	1
All	All	864/874 (99%)	648 (75%)	133 (15%)	83 (10%)	1	3

5 of 83 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	26	SER
1	L	51	ALA
1	L	168	SER
2	H	16	GLY
2	H	53	SER

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	K	176/183 (96%)	108 (61%)	68 (39%)	0	0
1	L	176/183 (96%)	125 (71%)	51 (29%)	0	2
2	H	185/190 (97%)	119 (64%)	66 (36%)	0	1
2	M	185/190 (97%)	132 (71%)	53 (29%)	0	2
All	All	722/746 (97%)	484 (67%)	238 (33%)	0	1

5 of 238 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	H	198	LEU
1	K	61	ARG
2	M	146	ASP
2	H	209	ASN
1	K	12	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such sidechains are listed below:

Mol	Chain	Res	Type
2	H	212	HIS
1	K	137	ASN
2	M	212	HIS
1	K	38	GLN
1	K	90	HIS

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 ⓘ

22 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
4	NAG	H	230	4	15,15,15	1.22	1 (6%)	21,21,21	2.65	9 (42%)
4	NAG	H	231	5,4	14,14,15	0.89	0	15,19,21	2.36	5 (33%)
5	BMA	H	232	3,4	11,11,12	1.17	1 (9%)	13,15,17	2.92	7 (53%)
3	MAN	H	233	3,5	11,11,12	1.49	3 (27%)	13,15,17	2.07	5 (38%)
3	MAN	H	234	3,5	11,11,12	0.94	0	13,15,17	2.37	4 (30%)
3	MAN	H	235	3	11,11,12	1.84	3 (27%)	13,15,17	3.68	9 (69%)
3	MAN	H	236	3	11,11,12	2.16	5 (45%)	13,15,17	2.90	4 (30%)
3	MAN	H	237	3	11,11,12	1.26	2 (18%)	13,15,17	4.32	9 (69%)
3	MAN	H	238	3,2	11,11,12	1.40	2 (18%)	13,15,17	3.42	6 (46%)
4	NAG	K	214	4	15,15,15	1.64	4 (26%)	21,21,21	2.71	9 (42%)
4	NAG	K	215	5,4	14,14,15	0.98	1 (7%)	15,19,21	2.80	7 (46%)
5	BMA	K	216	3,4	11,11,12	1.38	2 (18%)	13,15,17	2.73	8 (61%)
3	MAN	K	217	5	11,11,12	1.11	1 (9%)	13,15,17	2.67	5 (38%)
3	MAN	K	218	3,5	11,11,12	1.23	1 (9%)	13,15,17	3.52	7 (53%)
3	MAN	K	219	3	11,11,12	1.74	3 (27%)	13,15,17	2.32	6 (46%)
3	MAN	L	214	3	11,11,12	1.20	2 (18%)	13,15,17	2.91	7 (53%)
3	MAN	L	215	3	11,11,12	1.51	1 (9%)	13,15,17	2.33	4 (30%)
3	MAN	M	230	3,2	11,11,12	1.64	3 (27%)	13,15,17	3.54	7 (53%)
3	MAN	M	231	3	11,11,12	1.19	1 (9%)	13,15,17	3.02	6 (46%)
3	MAN	M	232	3	11,11,12	1.03	0	13,15,17	1.97	5 (38%)
3	MAN	M	233	3	11,11,12	1.40	1 (9%)	13,15,17	2.60	5 (38%)
3	MAN	M	234	-	11,11,12	0.91	0	13,15,17	3.30	5 (38%)

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
4	NAG	H	230	4	-	0/6/26/26	0/1/1/1
4	NAG	H	231	5,4	-	0/6/23/26	0/1/1/1
5	BMA	H	232	3,4	-	0/2/19/22	0/1/1/1
3	MAN	H	233	3,5	-	0/2/19/22	0/1/1/1
3	MAN	H	234	3,5	-	0/2/19/22	0/1/1/1
3	MAN	H	235	3	-	0/2/19/22	0/1/1/1
3	MAN	H	236	3	-	0/2/19/22	0/1/1/1
3	MAN	H	237	3	-	0/2/19/22	0/1/1/1
3	MAN	H	238	3,2	-	0/2/19/22	0/1/1/1
4	NAG	K	214	4	-	0/6/26/26	0/1/1/1
4	NAG	K	215	5,4	-	0/6/23/26	0/1/1/1
5	BMA	K	216	3,4	-	0/2/19/22	0/1/1/1
3	MAN	K	217	5	-	0/2/19/22	0/1/1/1
3	MAN	K	218	3,5	-	0/2/19/22	0/1/1/1
3	MAN	K	219	3	-	0/2/19/22	0/1/1/1
3	MAN	L	214	3	-	0/2/19/22	0/1/1/1
3	MAN	L	215	3	-	0/2/19/22	0/1/1/1
3	MAN	M	230	3,2	-	0/2/19/22	0/1/1/1
3	MAN	M	231	3	-	0/2/19/22	0/1/1/1
3	MAN	M	232	3	-	0/2/19/22	0/1/1/1
3	MAN	M	233	3	-	0/2/19/22	0/1/1/1
3	MAN	M	234	-	-	0/2/19/22	0/1/1/1

The worst 5 of 37 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	215	MAN	O5-C1	-4.36	1.36	1.43
3	M	230	MAN	C2-C3	-3.13	1.48	1.52
3	H	238	MAN	O4-C4	-3.04	1.35	1.43
3	M	230	MAN	O5-C1	-2.95	1.38	1.43
3	H	233	MAN	C4-C5	-2.75	1.47	1.53

The worst 5 of 139 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	238	MAN	C1-C2-C3	-8.78	98.51	109.65
3	M	230	MAN	O2-C2-C3	-8.19	94.08	110.17
3	M	231	MAN	C1-C2-C3	-7.35	100.33	109.65
3	H	237	MAN	O3-C3-C4	-6.97	95.20	110.36
3	L	214	MAN	C6-C5-C4	-6.12	98.69	113.00

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

16 monomers are involved in 41 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	H	232	BMA	4	0
3	H	233	MAN	2	0
3	H	234	MAN	11	0
3	H	235	MAN	3	0
3	H	236	MAN	2	0
3	H	237	MAN	7	0
3	H	238	MAN	4	0
5	K	216	BMA	3	0
3	K	217	MAN	4	0
3	K	218	MAN	8	0
3	K	219	MAN	3	0
3	L	214	MAN	6	0
3	M	230	MAN	8	0
3	M	231	MAN	5	0
3	M	232	MAN	5	0
3	M	234	MAN	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	H	1
2	M	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	M	112:SER	C	113:PRO	N	2.05
1	H	112:SER	C	113:PRO	N	1.70

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.