



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:45 AM GMT

PDB ID : 2E7Q
Title : Crystal structure of basic winged bean lectin in complex with b blood group trisaccharide
Authors : Kulkarni, K.A.; Katiyar, S.; Surolia, A.; Vijayan, M.; Suguna, K.
Deposited on : 2007-01-12
Resolution : 2.75 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 (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

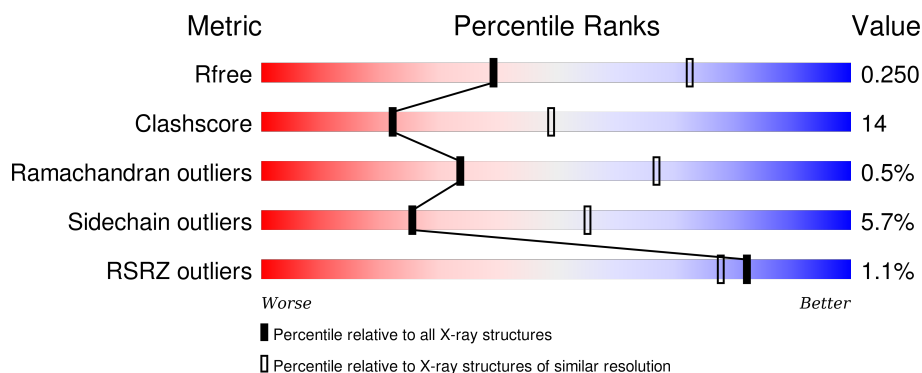
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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.

Mol	Chain	Length	Quality of chain
1	A	237	<div> <div>2%</div> <div>76%</div> <div>22%</div> <div>.</div> </div>
1	B	237	<div> <div>72%</div> <div>24%</div> <div>.</div> </div>
1	C	237	<div> <div>72%</div> <div>27%</div> <div>.</div> </div>
1	D	237	<div> <div>2%</div> <div>71%</div> <div>26%</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	NAG	C	501	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7908 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 Basic agglutinin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	0	0	0
			1839	1182	308	349			
1	B	237	Total	C	N	O	0	0	0
			1823	1172	303	348			
1	C	237	Total	C	N	O	0	0	0
			1817	1169	302	346			
1	D	237	Total	C	N	O	0	0	0
			1791	1154	295	342			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	3	Total	C	N	O	0	0
			38	22	2	14		
2	A	3	Total	C	N	O	0	0
			38	22	2	14		
2	B	3	Total	C	N	O	0	0
			38	22	2	14		
2	C	3	Total	C	N	O	0	0
			38	22	2	14		
2	D	3	Total	C	N	O	0	0
			38	22	2	14		
2	D	3	Total	C	N	O	0	0
			38	22	2	14		

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

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	3	Total	C	O	0	0
			34	18	16		
3	B	3	Total	C	O	0	0
			34	18	16		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	3	Total	C	O	0	0
			34	18	16		

- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		
6	A	1	Total	Ca	0	0
			1	1		
6	D	1	Total	Ca	0	0
			1	1		
6	C	1	Total	Ca	0	0
			1	1		

- Molecule 7 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Mn	0	0
			1	1		
7	A	1	Total	Mn	0	0
			1	1		
7	D	1	Total	Mn	0	0
			1	1		
7	C	1	Total	Mn	0	0
			1	1		

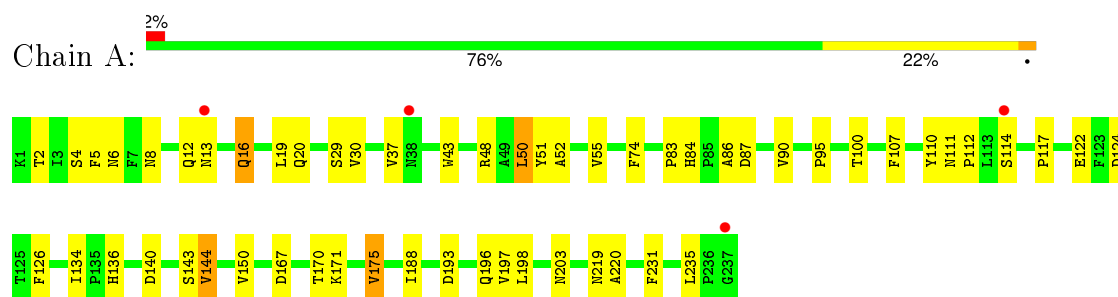
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	66	Total 66	O 66	0	0
8	B	52	Total 52	O 52	0	0
8	C	59	Total 59	O 59	0	0
8	D	57	Total 57	O 57	0	0

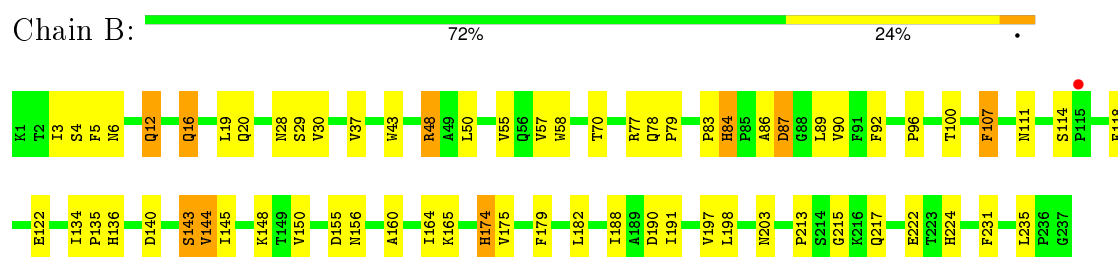
3 Residue-property plots [i](#)

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.

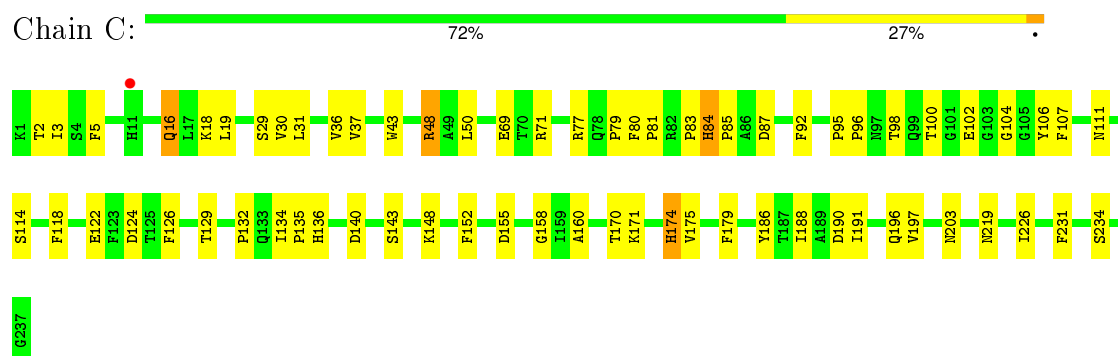
• Molecule 1: Basic agglutinin



• Molecule 1: Basic agglutinin

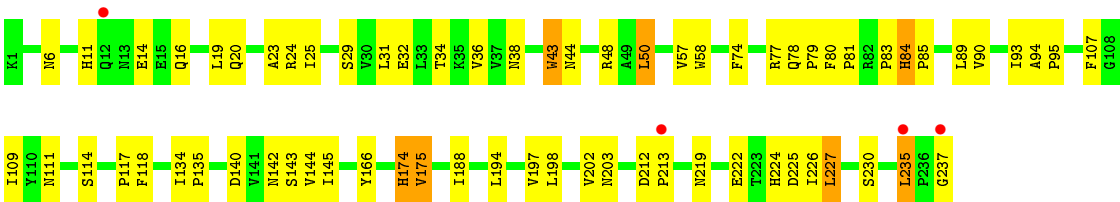


• Molecule 1: Basic agglutinin



• Molecule 1: Basic agglutinin





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	157.41Å 90.55Å 73.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.74 – 2.75 29.74 – 2.71	Depositor EDS
% Data completeness (in resolution range)	97.1 (29.74-2.75) 94.6 (29.74-2.71)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.72Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.198 , 0.258 0.192 , 0.250	Depositor DCC
R_{free} test set	1061 reflections (4.07%)	DCC
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.755	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 51.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 27583 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7908	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MN, NAG, GLA, CA, GLC, NDG, 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	0.43	0/1892	0.69	1/2589 (0.0%)
1	B	0.47	0/1876	0.70	1/2570 (0.0%)
1	C	0.44	0/1870	0.69	1/2562 (0.0%)
1	D	0.44	0/1843	0.70	1/2530 (0.0%)
All	All	0.44	0/7481	0.69	4/10251 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	29	SER	N-CA-C	5.69	126.36	111.00
1	A	29	SER	N-CA-C	5.54	125.96	111.00
1	C	29	SER	N-CA-C	5.14	124.89	111.00
1	D	29	SER	N-CA-C	5.09	124.73	111.00

There are no chirality outliers.

There are no planarity outliers.

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	1839	0	1785	42	0
1	B	1823	0	1750	45	0
1	C	1817	0	1740	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1791	0	1697	55	0
2	A	76	0	68	1	0
2	B	38	0	34	3	0
2	C	38	0	34	1	0
2	D	76	0	68	7	0
3	A	34	0	30	3	0
3	B	34	0	30	2	0
3	C	34	0	30	4	0
4	B	38	0	34	4	0
5	C	28	0	25	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
8	A	66	0	0	1	0
8	B	52	0	0	1	0
8	C	59	0	0	0	0
8	D	57	0	0	2	0
All	All	7908	0	7325	207	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 14.

All (207) 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)
3:A:401:GAL:C1	3:A:402:GLC:H62	2.04	0.86
1:C:16:GLN:HA	1:C:16:GLN:HE21	1.42	0.85
1:B:145:ILE:HB	1:C:196:GLN:HE22	1.44	0.82
1:D:197:VAL:HG23	1:D:198:LEU:HG	1.61	0.82
1:C:175:VAL:HG22	1:C:188:ILE:HG22	1.58	0.82
1:D:57:VAL:HG23	1:D:58:TRP:HD1	1.47	0.78
2:D:601:NAG:O4	2:D:602:FUC:H5	1.85	0.77
1:A:6:ASN:HD21	1:A:8:ASN:HD21	1.34	0.76
1:B:16:GLN:HE21	1:B:16:GLN:HA	1.52	0.75
1:C:155:ASP:OD2	1:C:158:GLY:HA3	1.87	0.75
3:C:401:GAL:C1	3:C:402:GLC:H61	2.16	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:VAL:HG22	1:A:188:ILE:HG22	1.69	0.74
1:A:16:GLN:HA	1:A:16:GLN:HE21	1.53	0.73
4:B:601:NAG:H4	4:B:603:NDG:H8C1	1.70	0.73
4:B:601:NAG:H62	4:B:603:NDG:C1	2.19	0.72
2:D:602:FUC:H62	2:D:603:NAG:H2	1.69	0.72
2:A:601:NAG:H62	2:A:603:NAG:H82	1.72	0.70
1:D:175:VAL:HG22	1:D:188:ILE:HG22	1.72	0.70
2:D:501:NAG:H61	2:D:503:NAG:H82	1.74	0.69
1:A:16:GLN:HE21	1:A:52:ALA:HB3	1.59	0.67
1:A:197:VAL:HG22	1:A:198:LEU:HG	1.77	0.67
3:B:401:GAL:C1	3:B:402:GLC:H61	2.26	0.66
1:B:48:ARG:HD2	1:B:100:THR:OG1	1.97	0.65
1:B:145:ILE:HB	1:C:196:GLN:NE2	2.10	0.65
1:B:213:PRO:HB3	4:B:601:NAG:H82	1.79	0.65
1:A:6:ASN:HD21	1:A:8:ASN:ND2	1.96	0.64
1:D:24:ARG:HH11	1:D:24:ARG:HG2	1.62	0.64
1:B:20:GLN:OE1	1:B:48:ARG:HD3	1.97	0.64
1:A:140:ASP:HB3	1:A:143:SER:O	1.98	0.63
1:C:48:ARG:HD2	1:C:100:THR:OG1	1.98	0.63
1:C:16:GLN:HA	1:C:16:GLN:NE2	2.12	0.63
2:D:602:FUC:H5	2:D:603:NAG:C1	2.30	0.62
1:B:78:GLN:HB2	1:B:156:ASN:HD21	1.64	0.61
1:D:140:ASP:HB3	1:D:143:SER:O	2.01	0.61
1:D:43:TRP:CZ3	1:D:213:PRO:HA	2.36	0.60
2:D:602:FUC:C6	2:D:603:NAG:H2	2.31	0.60
1:C:19:LEU:N	1:C:19:LEU:HD12	2.16	0.60
1:B:90:VAL:HG12	1:B:122:GLU:HA	1.84	0.60
1:B:150:VAL:HG12	1:B:188:ILE:HD12	1.83	0.60
1:D:38:ASN:N	1:D:38:ASN:HD22	2.00	0.59
1:B:111:ASN:ND2	1:B:114:SER:H	2.00	0.59
1:A:193:ASP:HB3	1:A:196:GLN:HB2	1.84	0.59
1:B:89:LEU:C	1:B:89:LEU:HD12	2.24	0.58
1:D:83:PRO:HD2	1:D:219:ASN:HB2	1.85	0.58
1:C:175:VAL:HG22	1:C:188:ILE:CG2	2.31	0.58
1:C:96:PRO:HA	1:C:203:ASN:OD1	2.03	0.58
1:A:16:GLN:NE2	1:A:52:ALA:HB3	2.20	0.57
1:A:20:GLN:OE1	1:A:48:ARG:HD3	2.04	0.57
1:A:175:VAL:HG22	1:A:188:ILE:CG2	2.34	0.56
1:C:126:PHE:CE2	3:C:400:GLA:H3	2.41	0.56
1:D:78:GLN:HG2	1:D:85:PRO:HB3	1.87	0.55
1:A:111:ASN:ND2	1:A:114:SER:HB3	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:GLN:HA	1:A:16:GLN:NE2	2.21	0.55
1:A:16:GLN:CA	1:A:16:GLN:HE21	2.18	0.55
1:C:140:ASP:HB3	1:C:143:SER:O	2.06	0.55
1:C:3:ILE:O	1:C:3:ILE:HG23	2.07	0.55
1:C:134:ILE:HD12	1:C:134:ILE:C	2.27	0.55
1:C:95:PRO:O	1:C:98:THR:HG23	2.07	0.55
1:C:50:LEU:HD11	1:C:92:PHE:CZ	2.42	0.54
1:C:102:GLU:HB3	1:C:106:TYR:HD2	1.72	0.54
1:D:117:PRO:HA	1:D:142:ASN:OD1	2.07	0.54
1:D:77:ARG:O	1:D:79:PRO:HD3	2.06	0.54
1:D:134:ILE:HD12	1:D:134:ILE:C	2.28	0.54
1:C:111:ASN:ND2	1:C:114:SER:H	2.06	0.54
1:C:134:ILE:O	1:C:134:ILE:HD12	2.08	0.53
1:B:28:ASN:HA	8:B:1307:HOH:O	2.07	0.53
1:C:16:GLN:CA	1:C:16:GLN:HE21	2.16	0.53
1:B:140:ASP:HB3	1:B:143:SER:O	2.08	0.53
1:B:197:VAL:HG23	1:B:198:LEU:HG	1.91	0.53
1:D:95:PRO:HD3	1:D:117:PRO:O	2.08	0.53
1:D:134:ILE:HD12	1:D:134:ILE:O	2.09	0.53
1:A:110:TYR:CE2	1:A:112:PRO:HG3	2.44	0.53
1:D:175:VAL:HG22	1:D:188:ILE:CG2	2.37	0.52
1:D:34:THR:HG21	1:D:224:HIS:ND1	2.24	0.52
1:D:203:ASN:HA	8:D:3330:HOH:O	2.10	0.52
1:D:77:ARG:NH1	1:D:79:PRO:HG3	2.24	0.52
2:B:501:NAG:H82	2:B:502:FUC:C1	2.39	0.52
2:C:502:FUC:H3	2:C:503:NAG:H61	1.92	0.52
2:D:501:NAG:C6	2:D:503:NAG:H82	2.37	0.52
1:D:16:GLN:HA	1:D:16:GLN:HE21	1.75	0.52
1:C:122:GLU:OE2	1:C:124:ASP:HB2	2.10	0.51
1:B:90:VAL:HG12	1:B:122:GLU:CB	2.39	0.51
1:D:212:ASP:OD1	1:D:213:PRO:HD2	2.11	0.51
1:D:83:PRO:HG2	1:D:84:HIS:CD2	2.46	0.51
1:A:144:VAL:O	1:A:144:VAL:HG13	2.10	0.51
1:B:5:PHE:CZ	1:B:231:PHE:HB3	2.46	0.51
1:C:31:LEU:HB3	1:C:226:ILE:HB	1.93	0.51
1:A:126:PHE:CE2	3:A:400:GLA:H3	2.46	0.51
1:D:84:HIS:CD2	1:D:84:HIS:O	2.64	0.50
1:D:235:LEU:HD22	1:D:237:GLY:H	1.76	0.50
1:D:77:ARG:HH11	1:D:79:PRO:HG3	1.76	0.50
1:D:20:GLN:OE1	1:D:48:ARG:NH1	2.41	0.50
1:B:70:THR:HG22	1:B:164:ILE:HB	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:501:NAG:H3	2:B:502:FUC:C5	2.41	0.49
1:D:227:LEU:CD1	1:D:227:LEU:N	2.75	0.49
1:D:23:ALA:O	1:D:24:ARG:HG2	2.12	0.49
1:D:32:GLU:HG2	1:D:225:ASP:OD1	2.11	0.49
1:A:37:VAL:HG23	1:A:37:VAL:O	2.12	0.49
1:A:203:ASN:HA	8:A:635:HOH:O	2.13	0.49
2:D:602:FUC:H61	2:D:603:NAG:H82	1.95	0.49
1:A:134:ILE:HD12	1:A:134:ILE:O	2.13	0.49
1:D:6:ASN:OD1	1:D:230:SER:HB3	2.13	0.48
1:D:16:GLN:HA	1:D:16:GLN:NE2	2.28	0.48
1:B:50:LEU:HD11	1:B:92:PHE:HZ	1.78	0.48
1:A:150:VAL:HG12	1:A:188:ILE:HD12	1.95	0.48
1:C:152:PHE:HA	1:C:186:TYR:CE1	2.49	0.48
1:D:222:GLU:HG3	1:D:224:HIS:CE1	2.49	0.48
1:B:134:ILE:HG22	1:B:136:HIS:CE1	2.49	0.48
1:A:16:GLN:HB3	1:A:51:TYR:CE1	2.49	0.47
1:B:84:HIS:CD2	1:B:84:HIS:O	2.67	0.47
4:B:601:NAG:C6	4:B:603:NDG:C1	2.92	0.47
1:D:24:ARG:HG2	1:D:24:ARG:NH1	2.27	0.47
1:B:222:GLU:OE2	1:B:224:HIS:NE2	2.46	0.47
1:B:90:VAL:HG12	1:B:122:GLU:CA	2.43	0.47
1:C:83:PRO:HG2	1:C:84:HIS:CD2	2.48	0.47
1:C:174:HIS:CD2	1:C:174:HIS:N	2.83	0.47
1:C:87:ASP:OD1	1:C:104:GLY:HA2	2.14	0.47
1:A:83:PRO:HD2	1:A:219:ASN:HB2	1.96	0.47
1:D:38:ASN:ND2	1:D:38:ASN:N	2.62	0.47
1:C:134:ILE:HB	1:C:135:PRO:HA	1.96	0.47
1:B:84:HIS:HD2	1:B:84:HIS:O	1.98	0.47
1:C:69:GLU:OE2	1:C:71:ARG:NE	2.48	0.47
1:C:175:VAL:CG2	1:C:188:ILE:HG22	2.38	0.47
1:A:150:VAL:HG11	1:A:188:ILE:HB	1.96	0.47
1:A:84:HIS:O	1:A:220:ALA:HA	2.15	0.47
1:B:215:GLY:O	1:B:217:GLN:HG2	2.14	0.47
1:D:50:LEU:CD1	1:D:203:ASN:HB3	2.45	0.47
1:C:132:PRO:HG2	1:C:136:HIS:ND1	2.30	0.47
1:B:20:GLN:OE1	1:B:48:ARG:CD	2.63	0.46
1:A:110:TYR:HE2	1:A:112:PRO:HG3	1.79	0.46
1:A:90:VAL:HG12	1:A:122:GLU:CB	2.46	0.46
1:D:174:HIS:HB2	8:D:3304:HOH:O	2.14	0.46
1:B:174:HIS:CD2	1:B:174:HIS:N	2.84	0.46
1:C:18:LYS:HE3	1:C:18:LYS:HB2	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:GLU:HG3	1:B:224:HIS:CE1	2.51	0.45
1:D:16:GLN:CA	1:D:16:GLN:HE21	2.28	0.45
1:B:134:ILE:HB	1:B:135:PRO:HA	1.97	0.45
1:A:111:ASN:HD22	1:A:114:SER:HB3	1.80	0.45
2:B:501:NAG:H3	2:B:502:FUC:H5	1.97	0.45
1:D:89:LEU:C	1:D:89:LEU:HD12	2.37	0.45
1:C:160:ALA:HB2	1:C:179:PHE:CE1	2.52	0.45
1:A:170:THR:O	1:A:171:LYS:HB2	2.17	0.45
1:A:197:VAL:CG2	1:A:198:LEU:HG	2.45	0.45
1:B:86:ALA:HA	1:B:87:ASP:HA	1.82	0.45
1:D:31:LEU:HB3	1:D:226:ILE:HB	1.99	0.45
1:D:166:TYR:CE2	1:D:194:LEU:HD12	2.52	0.45
1:B:50:LEU:HD11	1:B:92:PHE:CZ	2.52	0.44
1:C:2:THR:HG22	1:C:234:SER:OG	2.16	0.44
1:D:44:ASN:HB2	1:D:213:PRO:HG3	1.99	0.44
1:B:57:VAL:HG23	1:B:58:TRP:HD1	1.81	0.44
1:C:190:ASP:OD1	1:C:191:ILE:N	2.50	0.44
1:D:93:ILE:CG2	1:D:202:VAL:HG21	2.47	0.44
3:B:401:GAL:C1	3:B:402:GLC:C6	2.95	0.44
1:B:96:PRO:HA	1:B:203:ASN:OD1	2.17	0.44
1:C:36:VAL:O	1:C:36:VAL:HG13	2.16	0.44
1:C:77:ARG:CZ	1:C:79:PRO:HG3	2.48	0.44
1:C:124:ASP:HB3	1:C:136:HIS:CD2	2.52	0.44
1:A:74:PHE:CD1	1:A:74:PHE:C	2.90	0.44
1:D:57:VAL:HG23	1:D:58:TRP:CD1	2.39	0.44
1:C:77:ARG:O	1:C:79:PRO:HD3	2.18	0.43
1:B:77:ARG:O	1:B:79:PRO:HD3	2.17	0.43
1:C:37:VAL:O	1:C:37:VAL:HG23	2.18	0.43
1:A:50:LEU:N	1:A:50:LEU:HD22	2.34	0.43
1:A:84:HIS:CD2	1:A:126:PHE:HE1	2.37	0.43
1:D:212:ASP:OD1	1:D:213:PRO:CD	2.67	0.43
1:A:196:GLN:HE22	1:D:145:ILE:HB	1.83	0.43
1:D:74:PHE:CD1	1:D:74:PHE:C	2.91	0.43
1:D:90:VAL:HG21	1:D:109:ILE:HD13	2.00	0.43
1:D:118:PHE:C	1:D:118:PHE:CD1	2.91	0.42
1:C:5:PHE:CZ	1:C:231:PHE:HB3	2.53	0.42
1:A:134:ILE:HD12	1:A:134:ILE:C	2.40	0.42
1:C:148:LYS:HD3	1:C:190:ASP:OD2	2.20	0.42
1:B:118:PHE:C	1:B:118:PHE:CD1	2.93	0.42
1:D:11:HIS:O	1:D:14:GLU:HB2	2.19	0.42
3:C:401:GAL:O5	3:C:402:GLC:H61	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:ILE:O	1:B:4:SER:HB3	2.18	0.42
1:B:78:GLN:HB2	1:B:156:ASN:ND2	2.32	0.42
1:D:134:ILE:HB	1:D:135:PRO:HA	2.02	0.42
1:A:167:ASP:OD2	1:B:148:LYS:HE2	2.20	0.42
1:D:111:ASN:ND2	1:D:114:SER:OG	2.52	0.41
1:D:212:ASP:OD1	1:D:213:PRO:N	2.53	0.41
1:B:83:PRO:HG2	1:B:84:HIS:CD2	2.55	0.41
1:B:12:GLN:HE21	1:B:12:GLN:HB2	1.61	0.41
1:B:37:VAL:HG23	1:B:37:VAL:O	2.20	0.41
1:B:160:ALA:HB2	1:B:179:PHE:CE1	2.54	0.41
1:A:83:PRO:O	1:A:84:HIS:HB3	2.19	0.41
1:C:118:PHE:CD1	1:C:118:PHE:C	2.93	0.41
1:B:107:PHE:HE2	1:B:144:VAL:HG22	1.86	0.41
1:A:5:PHE:CZ	1:A:231:PHE:HB3	2.56	0.41
3:A:401:GAL:O5	3:A:402:GLC:H62	2.18	0.41
1:A:86:ALA:HA	1:A:87:ASP:HA	1.81	0.41
1:A:48:ARG:HD2	1:A:100:THR:OG1	2.21	0.41
1:B:50:LEU:HD22	1:B:50:LEU:N	2.35	0.41
1:C:83:PRO:HD2	1:C:219:ASN:HB2	2.03	0.41
1:A:124:ASP:HB3	1:A:136:HIS:CD2	2.55	0.41
1:B:190:ASP:OD1	1:B:191:ILE:N	2.53	0.41
1:A:95:PRO:HD3	1:A:117:PRO:O	2.20	0.41
1:C:170:THR:O	1:C:171:LYS:HB2	2.21	0.41
1:C:126:PHE:CZ	3:C:400:GLA:H5	2.56	0.40
1:B:155:ASP:HB2	1:B:182:LEU:HD11	2.03	0.40
1:D:94:ALA:HB1	1:D:95:PRO:CD	2.51	0.40
1:C:84:HIS:HA	1:C:85:PRO:HD3	1.93	0.40
1:D:235:LEU:CD2	1:D:237:GLY:H	2.33	0.40
1:C:84:HIS:CD2	1:C:84:HIS:O	2.75	0.40
1:C:80:PHE:HA	1:C:81:PRO:HD2	1.82	0.40
1:D:80:PHE:HA	1:D:81:PRO:HD2	1.94	0.40
1:D:19:LEU:HD21	1:D:25:ILE:HG13	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	235/237 (99%)	220 (94%)	13 (6%)	2 (1%)	21	52
1	B	235/237 (99%)	218 (93%)	16 (7%)	1 (0%)	39	72
1	C	235/237 (99%)	224 (95%)	10 (4%)	1 (0%)	39	72
1	D	235/237 (99%)	222 (94%)	12 (5%)	1 (0%)	39	72
All	All	940/948 (99%)	884 (94%)	51 (5%)	5 (0%)	34	67

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	ASN
1	A	107	PHE
1	C	107	PHE
1	D	107	PHE
1	B	107	PHE

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	204/206 (99%)	192 (94%)	12 (6%)	24	53
1	B	200/206 (97%)	184 (92%)	16 (8%)	15	37
1	C	198/206 (96%)	190 (96%)	8 (4%)	38	71
1	D	192/206 (93%)	183 (95%)	9 (5%)	32	64
All	All	794/824 (96%)	749 (94%)	45 (6%)	25	55

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR

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Mol	Chain	Res	Type
1	A	4	SER
1	A	12	GLN
1	A	16	GLN
1	A	19	LEU
1	A	30	VAL
1	A	43	TRP
1	A	50	LEU
1	A	55	VAL
1	A	144	VAL
1	A	175	VAL
1	A	235	LEU
1	B	6	ASN
1	B	12	GLN
1	B	16	GLN
1	B	19	LEU
1	B	30	VAL
1	B	43	TRP
1	B	48	ARG
1	B	55	VAL
1	B	84	HIS
1	B	87	ASP
1	B	143	SER
1	B	144	VAL
1	B	165	LYS
1	B	174	HIS
1	B	175	VAL
1	B	235	LEU
1	C	16	GLN
1	C	30	VAL
1	C	43	TRP
1	C	48	ARG
1	C	84	HIS
1	C	129	THR
1	C	174	HIS
1	C	197	VAL
1	D	36	VAL
1	D	43	TRP
1	D	50	LEU
1	D	84	HIS
1	D	144	VAL
1	D	174	HIS
1	D	175	VAL

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Mol	Chain	Res	Type
1	D	227	LEU
1	D	235	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	11	HIS
1	A	12	GLN
1	A	16	GLN
1	A	38	ASN
1	A	64	ASN
1	A	78	GLN
1	A	97	ASN
1	A	111	ASN
1	A	196	GLN
1	A	203	ASN
1	B	12	GLN
1	B	16	GLN
1	B	38	ASN
1	B	64	ASN
1	B	78	GLN
1	B	84	HIS
1	B	97	ASN
1	B	111	ASN
1	B	133	GLN
1	B	174	HIS
1	C	8	ASN
1	C	12	GLN
1	C	16	GLN
1	C	38	ASN
1	C	64	ASN
1	C	78	GLN
1	C	84	HIS
1	C	97	ASN
1	C	111	ASN
1	C	196	GLN
1	D	12	GLN
1	D	16	GLN
1	D	38	ASN
1	D	64	ASN
1	D	78	GLN

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Mol	Chain	Res	Type
1	D	84	HIS
1	D	111	ASN

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 ⓘ

32 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	$\# Z > 2$
3	GLA	A	400	3	11,11,12	0.67	0	14,15,17	0.77	1 (7%)
3	GAL	A	401	3	11,11,12	0.79	1 (9%)	14,15,17	0.62	0
3	GLC	A	402	3	12,12,12	0.63	0	17,17,17	0.49	0
2	NAG	A	501	1,2	14,14,15	0.55	0	15,19,21	0.71	0
2	FUC	A	502	2	10,10,11	0.70	0	14,14,16	0.62	0
2	NAG	A	503	2	14,14,15	0.48	0	15,19,21	0.87	1 (6%)
2	NAG	A	601	1,2	14,14,15	0.73	0	15,19,21	1.32	3 (20%)
2	FUC	A	602	2	10,10,11	0.74	0	14,14,16	0.93	1 (7%)
2	NAG	A	603	2	14,14,15	0.71	0	15,19,21	0.66	0
3	GLA	B	400	3	11,11,12	0.55	0	14,15,17	0.82	1 (7%)
3	GAL	B	401	3	11,11,12	0.67	0	14,15,17	0.67	0
3	GLC	B	402	3	12,12,12	0.49	0	17,17,17	0.43	0
2	NAG	B	501	1,2	14,14,15	0.69	0	15,19,21	0.86	0
2	FUC	B	502	2	10,10,11	0.61	0	14,14,16	0.60	0
2	NAG	B	503	2	14,14,15	0.70	0	15,19,21	0.79	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	B	601	1,4	14,14,15	0.72	0	15,19,21	0.94	1 (6%)
4	FUC	B	602	4	10,10,11	0.43	0	14,14,16	0.48	0
4	NDG	B	603	4	14,14,15	0.96	1 (7%)	15,19,21	0.92	1 (6%)
3	GLA	C	400	3	11,11,12	0.49	0	14,15,17	0.66	0
3	GAL	C	401	3	11,11,12	0.72	0	14,15,17	0.59	0
3	GLC	C	402	3	12,12,12	0.62	0	17,17,17	0.43	0
2	NAG	C	501	1,2	14,14,15	0.58	0	15,19,21	0.76	0
2	FUC	C	502	2	10,10,11	0.50	0	14,14,16	0.55	0
2	NAG	C	503	2	14,14,15	0.50	0	15,19,21	0.74	1 (6%)
5	NAG	C	601	1,5	14,14,15	0.60	0	15,19,21	0.92	1 (6%)
5	NAG	C	603	5	14,14,15	0.65	0	15,19,21	0.93	1 (6%)
2	NAG	D	501	1,2	14,14,15	0.62	0	15,19,21	0.64	0
2	FUC	D	502	2	10,10,11	0.62	0	14,14,16	0.61	0
2	NAG	D	503	2	14,14,15	0.71	0	15,19,21	0.71	0
2	NAG	D	601	1,2	14,14,15	0.58	0	15,19,21	0.92	1 (6%)
2	FUC	D	602	2	10,10,11	0.66	0	14,14,16	0.71	0
2	NAG	D	603	2	14,14,15	0.58	0	15,19,21	0.70	1 (6%)

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
3	GLA	A	400	3	-	0/2/19/22	0/1/1/1
3	GAL	A	401	3	-	0/2/19/22	0/1/1/1
3	GLC	A	402	3	-	0/2/22/22	0/1/1/1
2	NAG	A	501	1,2	-	0/6/23/26	0/1/1/1
2	FUC	A	502	2	-	0/0/17/20	0/1/1/1
2	NAG	A	503	2	-	0/6/23/26	0/1/1/1
2	NAG	A	601	1,2	-	0/6/23/26	0/1/1/1
2	FUC	A	602	2	-	0/0/17/20	0/1/1/1
2	NAG	A	603	2	-	0/6/23/26	0/1/1/1
3	GLA	B	400	3	-	0/2/19/22	0/1/1/1
3	GAL	B	401	3	-	0/2/19/22	0/1/1/1
3	GLC	B	402	3	-	0/2/22/22	0/1/1/1
2	NAG	B	501	1,2	-	0/6/23/26	0/1/1/1
2	FUC	B	502	2	-	0/0/17/20	0/1/1/1
2	NAG	B	503	2	-	0/6/23/26	0/1/1/1
4	NAG	B	601	1,4	-	0/6/23/26	0/1/1/1
4	FUC	B	602	4	-	0/0/17/20	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NDG	B	603	4	-	0/6/23/26	0/1/1/1
3	GLA	C	400	3	-	0/2/19/22	0/1/1/1
3	GAL	C	401	3	-	0/2/19/22	0/1/1/1
3	GLC	C	402	3	-	0/2/22/22	0/1/1/1
2	NAG	C	501	1,2	-	0/6/23/26	0/1/1/1
2	FUC	C	502	2	-	0/0/17/20	0/1/1/1
2	NAG	C	503	2	-	0/6/23/26	0/1/1/1
5	NAG	C	601	1,5	-	0/6/23/26	0/1/1/1
5	NAG	C	603	5	-	0/6/23/26	0/1/1/1
2	NAG	D	501	1,2	-	0/6/23/26	0/1/1/1
2	FUC	D	502	2	-	0/0/17/20	0/1/1/1
2	NAG	D	503	2	-	0/6/23/26	0/1/1/1
2	NAG	D	601	1,2	-	0/6/23/26	0/1/1/1
2	FUC	D	602	2	-	0/0/17/20	0/1/1/1
2	NAG	D	603	2	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	GAL	C2-C3	2.00	1.55	1.52
4	B	603	NDG	C1-C2	2.73	1.56	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	503	NAG	C2-N2-C7	-2.79	119.45	123.04
2	D	601	NAG	C2-N2-C7	-2.64	119.65	123.04
5	C	603	NAG	C2-N2-C7	-2.58	119.72	123.04
2	A	601	NAG	C2-N2-C7	-2.52	119.80	123.04
2	A	601	NAG	C4-C3-C2	-2.45	107.42	111.23
2	A	601	NAG	C3-C4-C5	-2.32	106.15	110.20
2	D	603	NAG	C2-N2-C7	-2.24	120.16	123.04
2	B	503	NAG	C2-N2-C7	-2.23	120.17	123.04
2	C	503	NAG	C2-N2-C7	-2.23	120.17	123.04
4	B	601	NAG	C2-N2-C7	-2.01	120.45	123.04
3	B	400	GLA	C1-O5-C5	2.03	114.82	112.25
5	C	601	NAG	C4-C3-C2	2.06	114.43	111.23
4	B	603	NDG	C4-C3-C2	2.18	114.61	111.23
3	A	400	GLA	C1-O5-C5	2.44	115.34	112.25
2	A	602	FUC	C1-C2-C3	2.93	113.01	109.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

21 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	400	GLA	1	0
3	A	401	GAL	2	0
3	A	402	GLC	2	0
2	A	601	NAG	1	0
2	A	603	NAG	1	0
3	B	401	GAL	2	0
3	B	402	GLC	2	0
2	B	501	NAG	3	0
2	B	502	FUC	3	0
4	B	601	NAG	4	0
4	B	603	NDG	3	0
3	C	400	GLA	2	0
3	C	401	GAL	2	0
3	C	402	GLC	2	0
2	C	502	FUC	1	0
2	C	503	NAG	1	0
2	D	501	NAG	2	0
2	D	503	NAG	2	0
2	D	601	NAG	1	0
2	D	602	FUC	5	0
2	D	603	NAG	4	0

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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

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, 95th 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	237/237 (100%)	-0.17	4 (1%) 73 68	14, 27, 48, 58	0
1	B	237/237 (100%)	-0.25	1 (0%) 93 92	13, 27, 49, 54	0
1	C	237/237 (100%)	-0.20	1 (0%) 93 92	19, 29, 52, 68	0
1	D	237/237 (100%)	-0.18	4 (1%) 73 68	17, 30, 54, 65	0
All	All	948/948 (100%)	-0.20	10 (1%) 82 78	13, 29, 51, 68	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	237	GLY	3.8
1	D	237	GLY	3.0
1	D	12	GLN	2.6
1	D	213	PRO	2.3
1	B	115	PRO	2.2
1	A	13	ASN	2.2
1	D	235	LEU	2.1
1	A	38	ASN	2.1
1	C	11	HIS	2.1
1	A	114	SER	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](#)

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, 95th 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	NAG	C	501	14/15	0.90	0.24	2.27	42,47,49,52	0
2	NAG	A	601	14/15	0.92	0.23	1.67	51,55,60,64	0
4	NAG	B	601	14/15	0.86	0.27	1.48	51,54,62,63	0
2	NAG	D	501	14/15	0.73	0.33	1.32	71,73,77,77	0
2	NAG	B	501	14/15	0.89	0.23	1.05	54,55,61,64	0
3	GLA	B	400	11/12	0.94	0.17	0.18	43,47,49,50	0
3	GLA	C	400	11/12	0.94	0.18	-0.12	54,55,57,58	0
3	GLA	A	400	11/12	0.87	0.16	-0.30	43,46,47,49	0
2	NAG	A	501	14/15	0.93	0.13	-0.93	41,45,49,51	0
2	FUC	C	502	10/11	0.91	0.27	-	54,56,57,59	0
4	NDG	B	603	14/15	0.72	0.36	-	65,67,69,69	0
3	GLC	C	402	12/12	0.87	0.25	-	68,72,72,73	0
5	NAG	C	601	14/15	0.82	0.23	-	59,62,69,69	0
3	GAL	B	401	11/12	0.78	0.23	-	52,56,58,58	0
2	NAG	D	601	14/15	0.90	0.29	-	52,57,60,65	0
2	FUC	A	502	10/11	0.93	0.16	-	50,52,52,53	0
2	NAG	A	503	14/15	0.89	0.23	-	56,58,61,63	0
2	FUC	A	602	10/11	0.86	0.29	-	61,62,63,63	0
3	GAL	A	401	11/12	0.86	0.20	-	54,62,64,64	0
2	FUC	B	502	10/11	0.79	0.29	-	68,70,72,73	0
2	NAG	B	503	14/15	0.85	0.30	-	63,65,67,68	0
2	NAG	C	503	14/15	0.92	0.18	-	46,48,52,55	0
2	NAG	D	503	14/15	0.78	0.33	-	79,80,81,81	0
2	NAG	D	603	14/15	0.83	0.30	-	69,72,74,75	0
4	FUC	B	602	10/11	0.94	0.22	-	47,50,51,51	0
3	GLC	A	402	12/12	0.77	0.26	-	65,68,69,70	0
2	FUC	D	602	10/11	0.83	0.34	-	61,63,63,64	0
2	NAG	A	603	14/15	0.72	0.38	-	69,71,73,75	0
3	GAL	C	401	11/12	0.88	0.19	-	60,64,64,66	0
3	GLC	B	402	12/12	0.79	0.23	-	60,61,63,63	0
5	NAG	C	603	14/15	0.71	0.33	-	70,72,73,73	0
2	FUC	D	502	10/11	0.67	0.40	-	79,80,80,81	0

6.4 Ligands ⓘ

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, 95th 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(\AA^2)	Q<0.9
6	CA	A	303	1/1	0.97	0.13	-0.34	26,26,26,26	0
6	CA	D	3303	1/1	0.99	0.11	-1.80	21,21,21,21	0
7	MN	C	2300	1/1	0.97	0.08	-2.01	38,38,38,38	0
6	CA	C	2303	1/1	0.99	0.10	-2.01	22,22,22,22	0
6	CA	B	1303	1/1	0.95	0.09	-2.43	19,19,19,19	0
7	MN	A	300	1/1	0.97	0.05	-2.68	29,29,29,29	0
7	MN	D	3300	1/1	0.98	0.08	-3.59	42,42,42,42	0
7	MN	B	1300	1/1	0.98	0.05	-3.82	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.