



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:03 PM GMT

PDB ID : 4GZT
Title : N2 neuraminidase D151G mutant of A/Tanzania/205/2010 H3N2 in complex with oseltamivir carboxylate
Authors : Zhu, X.; Wilson, I.A.
Deposited on : 2012-09-06
Resolution : 2.19 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

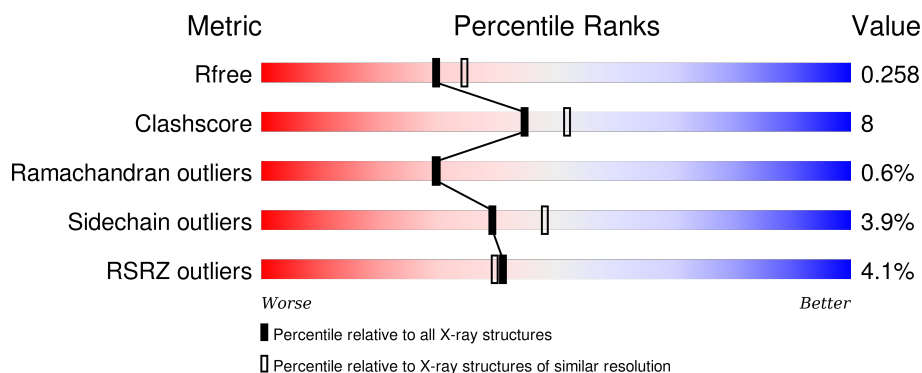
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (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.

Mol	Chain	Length	Quality of chain
1	A	393	<div> <div>3%</div> <div>76%</div> <div>21%</div> <div>..</div> </div>
1	B	393	<div> <div>5%</div> <div>79%</div> <div>18%</div> <div>...</div> </div>
1	C	393	<div> <div>4%</div> <div>77%</div> <div>20%</div> <div>..</div> </div>
1	D	393	<div> <div>4%</div> <div>80%</div> <div>18%</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
3	NAG	D	502	-	-	-	X

2 Entry composition [i](#)

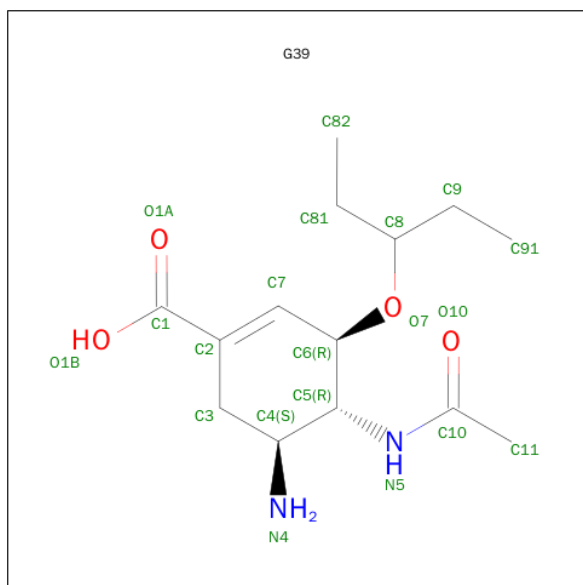
There are 9 unique types of molecules in this entry. The entry contains 13132 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 neuraminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	0	0	0
			2998	1858	531	587	22			
1	B	388	Total	C	N	O	S	0	0	0
			2998	1858	531	587	22			
1	C	388	Total	C	N	O	S	0	0	0
			2998	1858	531	587	22			
1	D	388	Total	C	N	O	S	0	0	0
			2998	1858	531	587	22			

- Molecule 2 is (3R,4R,5S)-4-(ACETYLAMINO)-5-AMINO-3-(PENTAN-3-YLOXY)CYCLOHEX-1-ENE-1-CARBOXYLIC ACID (three-letter code: G39) (formula: C₁₄H₂₄N₂O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			20	14	2	4		
2	B	1	Total	C	N	O	0	0
			20	14	2	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			20	14	2	4		
2	D	1	Total	C	N	O	0	0
			20	14	2	4		

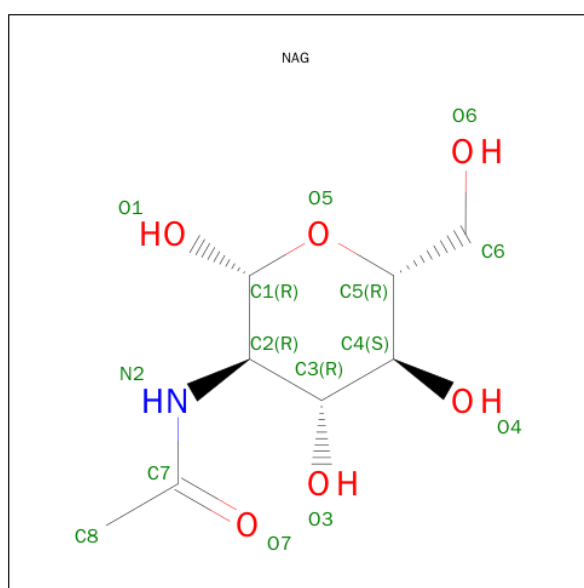
- Molecule 3 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	5	Total	C	N	O	0	0
			61	34	2	25		
3	B	5	Total	C	N	O	0	0
			61	34	2	25		
3	C	5	Total	C	N	O	0	0
			61	34	2	25		
3	D	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

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

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		

- 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	2	Total	Ca	0	0
			2	2		
6	D	1	Total	Ca	0	0
			1	1		
6	C	1	Total	Ca	0	0
			1	1		

- Molecule 7 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	2	Total	C	N	O	0	0
			24	14	1	9		

- Molecule 8 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	D	4	Total	C	N	O	0	0
			48	28	2	18		

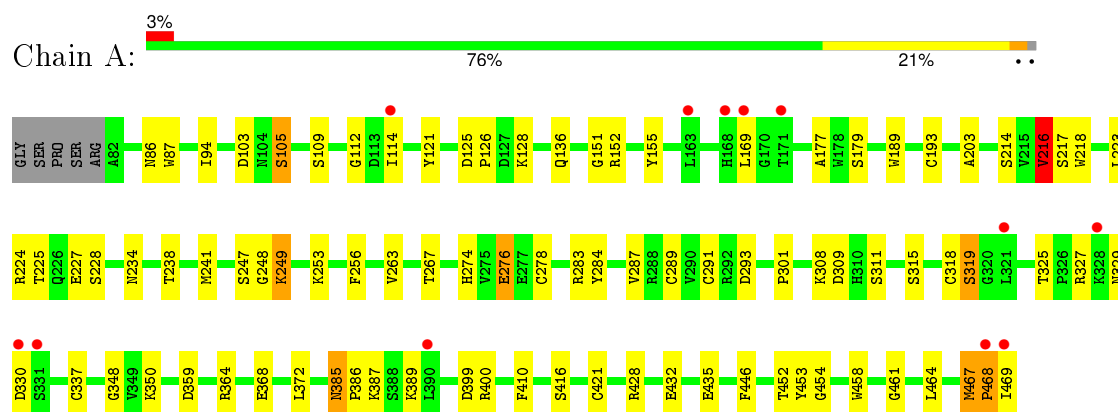
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	148	Total 148	O 148	0	0
9	B	147	Total 147	O 147	0	0
9	C	158	Total 158	O 158	0	0
9	D	146	Total 146	O 146	0	0

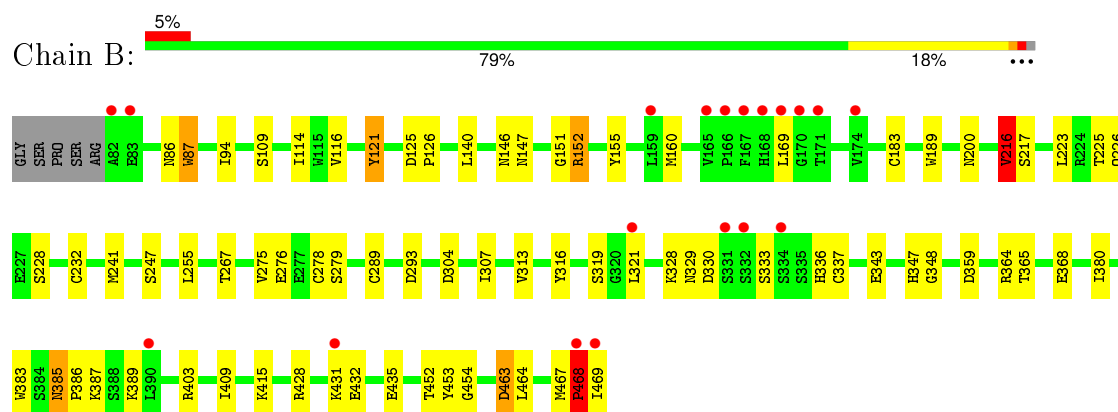
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

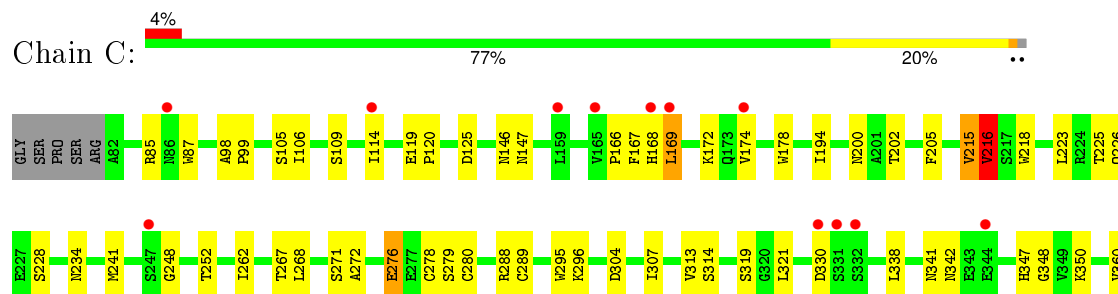
- Molecule 1: neuraminidase

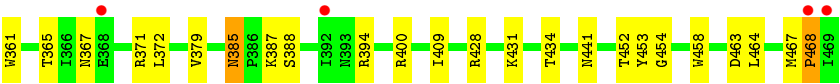


- Molecule 1: neuraminidase

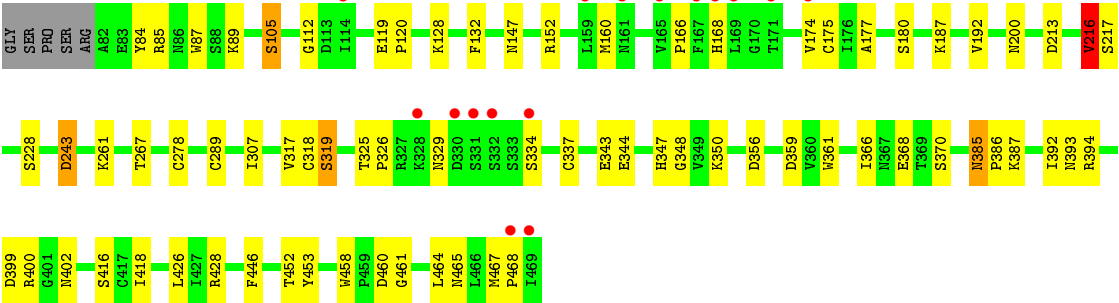
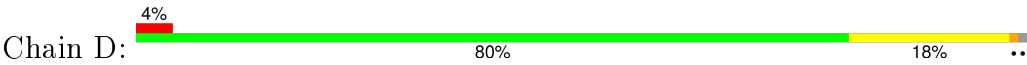


- Molecule 1: neuraminidase





● Molecule 1: neuraminidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.24Å 110.19Å 109.85Å 90.00° 98.27° 90.00°	Depositor
Resolution (Å)	49.14 – 2.19 49.14 – 2.19	Depositor EDS
% Data completeness (in resolution range)	88.9 (49.14-2.19) 88.9 (49.14-2.19)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.190 , 0.260 0.189 , 0.258	Depositor DCC
R_{free} test set	4467 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	1.527	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 88107 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13132	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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: G39, CA, BMA, NAG, 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	0.93	2/3065 (0.1%)	0.92	2/4156 (0.0%)
1	B	0.95	4/3065 (0.1%)	0.93	5/4156 (0.1%)
1	C	0.96	6/3065 (0.2%)	0.95	5/4156 (0.1%)
1	D	0.96	2/3065 (0.1%)	0.97	3/4156 (0.1%)
All	All	0.95	14/12260 (0.1%)	0.94	15/16624 (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	D	0	1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	218	TRP	CD2-CE2	5.88	1.48	1.41
1	D	87	TRP	CD2-CE2	5.80	1.48	1.41
1	B	87	TRP	CD2-CE2	5.78	1.48	1.41
1	C	87	TRP	CD2-CE2	5.68	1.48	1.41
1	B	383	TRP	CD2-CE2	5.65	1.48	1.41
1	C	178	TRP	CD2-CE2	5.59	1.48	1.41
1	D	361	TRP	CD2-CE2	5.49	1.48	1.41
1	B	121	TYR	CE1-CZ	5.49	1.45	1.38
1	C	458	TRP	CD2-CE2	5.46	1.48	1.41
1	C	167	PHE	CE2-CZ	5.29	1.47	1.37
1	A	87	TRP	CD2-CE2	5.27	1.47	1.41
1	C	361	TRP	CD2-CE2	5.24	1.47	1.41
1	A	189	TRP	CD2-CE2	5.13	1.47	1.41
1	B	189	TRP	CD2-CE2	5.00	1.47	1.41

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	216	VAL	CB-CA-C	-8.53	95.20	111.40
1	B	152	ARG	NE-CZ-NH1	8.35	124.47	120.30
1	A	216	VAL	CB-CA-C	-8.34	95.56	111.40
1	D	216	VAL	CB-CA-C	-8.12	95.97	111.40
1	B	216	VAL	CB-CA-C	-7.54	97.06	111.40
1	C	288	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	D	243	ASP	CB-CG-OD2	-6.15	112.76	118.30
1	B	152	ARG	NE-CZ-NH2	-5.72	117.44	120.30
1	C	169	LEU	CB-CG-CD1	-5.65	101.39	111.00
1	A	103	ASP	CB-CG-OD1	-5.56	113.30	118.30
1	C	216	VAL	CG1-CB-CG2	5.32	119.41	110.90
1	B	160	MET	CG-SD-CE	5.32	108.71	100.20
1	D	394	ARG	NE-CZ-NH1	5.21	122.90	120.30
1	B	403	ARG	NE-CZ-NH1	5.16	122.88	120.30
1	C	125	ASP	CB-CG-OD1	5.05	122.85	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	177	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	A	2998	0	2855	63	0
1	B	2998	0	2855	49	0
1	C	2998	0	2856	52	0
1	D	2998	0	2855	47	0
2	A	20	0	23	3	0
2	B	20	0	23	2	0
2	C	20	0	23	0	0
2	D	20	0	23	1	0
3	A	61	0	52	0	0
3	B	61	0	52	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	61	0	52	0	0
3	D	61	0	52	1	0
4	A	28	0	25	0	0
5	A	28	0	26	1	0
5	B	42	0	39	0	0
5	C	14	0	13	1	0
5	D	28	0	26	1	0
6	A	2	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	C	24	0	22	1	0
8	D	48	0	43	1	0
9	A	148	0	0	7	0
9	B	147	0	0	4	0
9	C	158	0	0	3	0
9	D	146	0	0	4	0
All	All	13132	0	11915	193	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 8.

All (193) 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:A:223:LEU:HD11	1:A:241:MET:HE1	1.52	0.91
1:C:109:SER:HB3	1:C:114:ILE:HB	1.52	0.90
1:A:452:THR:HB	1:B:216:VAL:HG22	1.56	0.88
1:C:85:ARG:HD3	9:C:621:HOH:O	1.75	0.84
1:A:216:VAL:HG22	1:D:452:THR:HB	1.58	0.83
1:B:467:MET:O	1:B:469:ILE:HG13	1.85	0.76
1:D:385:ASN:ND2	1:D:387:LYS:H	1.85	0.74
1:A:309:ASP:OD1	1:A:311:SER:OG	2.07	0.72
1:A:152:ARG:NH1	2:A:501:G39:O10	2.23	0.72
1:B:452:THR:HB	1:C:216:VAL:HG22	1.71	0.71
1:A:216:VAL:HG13	1:D:453:TYR:C	2.12	0.69
1:B:330:ASP:O	1:B:389:LYS:NZ	2.24	0.69
1:A:152:ARG:HD3	2:A:501:G39:H111	1.75	0.68
1:A:385:ASN:ND2	1:A:387:LYS:H	1.92	0.67
1:C:279:SER:HB3	1:C:409:ILE:HG22	1.78	0.65
1:A:467:MET:HB3	1:A:468:PRO:HD2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:146:ASN:O	1:C:147:ASN:HB2	1.95	0.65
3:D:506:BMA:H3	9:D:614:HOH:O	1.97	0.65
1:B:464:LEU:HA	1:B:467:MET:HG3	1.78	0.65
1:C:347:HIS:CD2	1:C:348:GLY:H	2.15	0.65
1:C:271:SER:HB3	1:C:338:LEU:O	1.97	0.64
1:B:152:ARG:NH1	2:B:501:G39:O10	2.30	0.64
1:A:330:ASP:O	1:A:389:LYS:NZ	2.25	0.63
1:C:147:ASN:ND2	7:C:508:FUC:H5	2.14	0.63
1:C:452:THR:HB	1:D:216:VAL:HG22	1.79	0.62
1:C:223:LEU:HD11	1:C:241:MET:HE2	1.81	0.62
1:A:225:THR:HB	1:A:241:MET:HG2	1.80	0.62
1:A:223:LEU:CD1	1:A:241:MET:HE1	2.26	0.62
1:B:152:ARG:HD3	2:B:501:G39:H111	1.80	0.61
1:D:385:ASN:HD22	1:D:387:LYS:H	1.48	0.61
1:B:321:LEU:HD22	1:B:330:ASP:OD1	1.99	0.61
1:C:304:ASP:HB2	1:C:313:VAL:CG2	2.30	0.61
1:C:347:HIS:CG	1:C:348:GLY:H	2.19	0.60
1:D:385:ASN:C	1:D:385:ASN:HD22	2.04	0.60
1:A:109:SER:HB3	1:A:114:ILE:HB	1.84	0.60
1:A:385:ASN:HD22	1:A:385:ASN:C	2.05	0.59
1:C:454:GLY:HA3	1:D:200:ASN:O	2.02	0.59
1:A:112:GLY:HA3	1:B:169:LEU:HD11	1.83	0.59
1:A:432:GLU:OE1	1:A:432:GLU:N	2.33	0.58
1:C:453:TYR:C	1:D:216:VAL:HG13	2.24	0.58
1:B:279:SER:HB3	1:B:409:ILE:HG22	1.86	0.58
1:B:109:SER:HB3	1:B:114:ILE:HB	1.85	0.58
1:B:385:ASN:ND2	1:B:387:LYS:H	2.00	0.58
1:A:216:VAL:O	9:A:623:HOH:O	2.17	0.58
1:D:278:CYS:HB3	1:D:289:CYS:HB3	1.86	0.58
1:A:216:VAL:HG22	1:D:452:THR:CB	2.31	0.58
1:C:278:CYS:HB3	1:C:289:CYS:HB3	1.86	0.57
1:B:385:ASN:HD21	1:B:387:LYS:HB2	1.69	0.57
1:A:464:LEU:HA	1:A:467:MET:HG3	1.86	0.57
1:A:325:THR:O	1:A:348:GLY:HA2	2.05	0.56
1:D:105:SER:HB2	9:D:744:HOH:O	2.04	0.56
1:D:147:ASN:ND2	8:D:510:FUC:H5	2.21	0.56
1:A:452:THR:HB	1:B:216:VAL:CG2	2.33	0.56
1:A:274:HIS:O	1:A:293:ASP:HA	2.05	0.56
1:B:347:HIS:CG	1:B:348:GLY:H	2.23	0.56
1:D:84:TYR:CE1	1:D:187:LYS:HD2	2.40	0.56
1:A:216:VAL:HG13	1:D:453:TYR:O	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:428:ARG:NH2	1:C:464:LEU:HG	2.22	0.55
1:A:256:PHE:O	1:A:263:VAL:HG22	2.07	0.55
1:C:205:PHE:CE1	1:C:262:ILE:HD11	2.42	0.54
1:B:146:ASN:O	1:B:147:ASN:HB2	2.07	0.54
1:A:177:ALA:HB2	1:A:193:CYS:HB3	1.88	0.54
1:B:151:GLY:HA3	9:B:746:HOH:O	2.06	0.54
1:D:85:ARG:HD3	9:D:665:HOH:O	2.07	0.54
1:C:365:THR:HG21	1:C:371:ARG:HA	1.91	0.53
1:D:465:ASN:H	1:D:465:ASN:HD22	1.56	0.53
1:B:247:SER:HB3	9:B:623:HOH:O	2.09	0.53
1:C:120:PRO:HG3	1:C:441:ASN:ND2	2.24	0.52
1:C:385:ASN:HD22	1:C:387:LYS:H	1.58	0.52
1:B:304:ASP:HB2	1:B:313:VAL:HG22	1.92	0.52
1:C:385:ASN:ND2	1:C:387:LYS:H	2.08	0.52
1:A:453:TYR:C	1:B:216:VAL:HG13	2.30	0.51
1:A:217:SER:HB2	9:A:602:HOH:O	2.09	0.51
1:A:203:ALA:O	1:A:214:SER:HA	2.11	0.51
1:B:328:LYS:HD2	1:B:343:GLU:OE1	2.11	0.51
1:D:356:ASP:O	1:D:359:ASP:HB2	2.10	0.51
1:C:225:THR:OG1	1:C:226:GLN:N	2.44	0.51
1:B:467:MET:HE3	1:B:468:PRO:HD2	1.93	0.50
1:B:223:LEU:HD11	1:B:241:MET:HE2	1.94	0.50
1:A:248:GLY:O	1:A:249:LYS:C	2.48	0.50
1:C:304:ASP:HB2	1:C:313:VAL:HG23	1.92	0.50
1:C:385:ASN:HD21	1:C:387:LYS:HB2	1.76	0.49
1:D:366:ILE:HG21	1:D:400:ARG:HB3	1.93	0.49
1:A:308:LYS:HB2	9:A:676:HOH:O	2.11	0.49
1:D:337:CYS:SG	1:D:386:PRO:HD3	2.52	0.49
1:D:343:GLU:O	1:D:344:GLU:HB2	2.12	0.49
1:D:399:ASP:CG	1:D:402:ASN:HD22	2.16	0.48
1:C:319:SER:OG	1:C:321:LEU:O	2.25	0.48
1:A:385:ASN:ND2	1:A:385:ASN:C	2.65	0.48
1:D:428:ARG:NH2	1:D:464:LEU:HG	2.28	0.48
1:A:283:ARG:O	1:A:284:TYR:C	2.52	0.48
1:D:326:PRO:HA	1:D:368:GLU:O	2.14	0.48
1:D:446:PHE:HZ	1:D:458:TRP:CE3	2.32	0.48
1:B:329:ASN:O	1:B:333:SER:HB3	2.14	0.48
1:C:248:GLY:HA2	1:C:295:TRP:CE2	2.49	0.47
1:B:463:ASP:O	1:B:467:MET:HG2	2.14	0.47
1:C:367:ASN:OD1	5:C:509:NAG:C7	2.62	0.47
1:A:224:ARG:NH2	1:A:276:GLU:OE2	2.37	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:THR:CG2	1:A:287:VAL:HG21	2.46	0.46
1:A:454:GLY:HA3	1:B:200:ASN:O	2.15	0.46
1:C:276:GLU:HG3	9:C:632:HOH:O	2.14	0.46
1:C:252:THR:HB	1:C:268:LEU:HD22	1.96	0.46
1:B:337:CYS:SG	1:B:386:PRO:HD3	2.56	0.46
1:B:364:ARG:HG3	1:B:365:THR:O	2.15	0.46
1:B:275:VAL:C	1:B:276:GLU:HG2	2.36	0.46
1:A:318:CYS:O	1:A:319:SER:C	2.54	0.46
1:B:121:TYR:CG	1:B:228:SER:HA	2.51	0.46
1:C:215:VAL:HG22	1:C:262:ILE:CD1	2.46	0.46
1:A:385:ASN:HD22	1:A:386:PRO:N	2.14	0.45
1:D:132:PHE:CE2	1:D:160:MET:HG3	2.50	0.45
1:A:136:GLN:NE2	9:A:619:HOH:O	2.45	0.45
1:C:347:HIS:CG	1:C:348:GLY:N	2.85	0.45
1:A:86:ASN:ND2	1:A:234:ASN:OD1	2.50	0.45
1:A:276:GLU:OE1	2:A:501:G39:C91	2.65	0.45
1:A:217:SER:CB	9:A:602:HOH:O	2.65	0.45
1:A:278:CYS:HB3	1:A:289:CYS:HB3	1.98	0.45
1:B:454:GLY:HA3	1:C:200:ASN:O	2.17	0.45
1:B:428:ARG:NH2	1:B:464:LEU:HG	2.32	0.45
1:A:315:SER:HB2	1:A:337:CYS:O	2.16	0.45
1:A:151:GLY:HA3	9:A:743:HOH:O	2.16	0.45
1:A:428:ARG:NH2	1:A:464:LEU:HG	2.32	0.45
1:B:432:GLU:N	1:B:432:GLU:OE1	2.51	0.44
1:B:183:CYS:SG	1:B:232:CYS:SG	3.15	0.44
1:D:89:LYS:HB2	1:D:418:ILE:HD11	1.99	0.44
1:A:400:ARG:NH1	5:A:510:NAG:O7	2.51	0.44
1:C:394:ARG:HD3	9:C:633:HOH:O	2.17	0.44
1:B:359:ASP:OD1	1:B:380:ILE:HA	2.17	0.44
1:C:304:ASP:HB2	1:C:313:VAL:HG22	1.99	0.44
1:D:325:THR:O	1:D:348:GLY:HA2	2.16	0.44
1:D:426:LEU:HD13	1:D:460:ASP:N	2.33	0.44
1:B:86:ASN:O	1:B:87:TRP:HB2	2.17	0.44
1:D:318:CYS:O	1:D:319:SER:C	2.57	0.43
1:B:453:TYR:C	1:C:216:VAL:HG13	2.37	0.43
1:C:272:ALA:HB2	1:C:314:SER:HB2	1.99	0.43
1:D:166:PRO:O	1:D:168:HIS:HD2	2.01	0.43
1:C:360:VAL:HG12	1:C:379:VAL:HB	2.00	0.43
1:B:321:LEU:CD2	1:B:330:ASP:OD1	2.64	0.43
1:B:125:ASP:HB2	1:B:126:PRO:CD	2.48	0.43
1:B:116:VAL:CG2	1:B:140:LEU:HA	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:400:ARG:NH1	5:D:512:NAG:O7	2.52	0.43
1:A:410:PHE:CZ	1:A:421:CYS:HB2	2.53	0.43
1:B:225:THR:OG1	1:B:226:GLN:N	2.50	0.43
1:D:119:GLU:N	1:D:120:PRO:CD	2.82	0.43
1:A:105:SER:HB2	9:A:604:HOH:O	2.19	0.43
1:D:213:ASP:OD2	1:D:261:LYS:HD2	2.18	0.43
1:C:453:TYR:O	1:D:216:VAL:HG13	2.19	0.42
1:C:194:ILE:HD12	1:C:225:THR:HG22	2.01	0.42
1:A:276:GLU:O	1:A:291:CYS:HB3	2.20	0.42
1:D:174:VAL:O	1:D:175:CYS:HB3	2.20	0.42
1:D:228:SER:HB3	1:D:350:LYS:HE2	2.00	0.42
1:B:217:SER:HB2	9:B:611:HOH:O	2.17	0.42
1:D:329:ASN:HA	1:D:368:GLU:HG2	2.00	0.42
1:D:119:GLU:N	1:D:120:PRO:HD3	2.34	0.42
1:A:223:LEU:HD11	1:A:241:MET:CE	2.37	0.42
1:D:217:SER:OG	1:D:243:ASP:OD2	2.28	0.42
1:A:121:TYR:CG	1:A:228:SER:HA	2.54	0.42
1:D:317:VAL:HA	9:D:692:HOH:O	2.19	0.42
1:D:464:LEU:HA	1:D:467:MET:HG3	2.01	0.41
1:C:172:LYS:HD3	1:C:174:VAL:HG12	2.02	0.41
1:C:394:ARG:HH11	1:C:394:ARG:HD3	1.73	0.41
1:A:169:LEU:HD11	1:D:112:GLY:HA3	2.02	0.41
1:C:341:ASN:O	1:C:342:ASN:HB2	2.20	0.41
1:A:179:SER:OG	1:A:227:GLU:OE2	2.30	0.41
1:B:468:PRO:HD3	9:B:709:HOH:O	2.19	0.41
1:C:280:CYS:HA	1:C:289:CYS:HA	2.01	0.41
1:C:169:LEU:HD23	1:C:169:LEU:HA	1.94	0.41
1:A:125:ASP:HB2	1:A:126:PRO:CD	2.51	0.41
1:D:180:SER:HA	1:D:192:VAL:O	2.20	0.41
1:A:446:PHE:HZ	1:A:458:TRP:CE3	2.38	0.41
1:B:385:ASN:HD22	1:B:385:ASN:C	2.23	0.41
1:C:98:ALA:HA	1:C:99:PRO:HD3	1.93	0.41
1:D:347:HIS:CG	1:D:348:GLY:H	2.39	0.41
1:A:169:LEU:HA	1:A:169:LEU:HD23	1.90	0.41
1:C:228:SER:HB3	1:C:350:LYS:HE2	2.03	0.41
1:B:278:CYS:HB3	1:B:289:CYS:HB3	2.02	0.41
1:A:218:TRP:CE2	1:A:253:LYS:HD2	2.55	0.41
1:B:255:LEU:HD12	1:B:255:LEU:N	2.35	0.41
1:D:152:ARG:HD3	2:D:501:G39:H111	2.03	0.41
1:C:119:GLU:N	1:C:120:PRO:HD3	2.36	0.40
1:C:385:ASN:HD22	1:C:385:ASN:C	2.24	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:296:LYS:O	1:C:342:ASN:HA	2.20	0.40
1:C:106:ILE:HG23	1:C:106:ILE:HD12	1.76	0.40
1:D:392:ILE:O	1:D:393:ASN:HB2	2.21	0.40
1:A:293:ASP:HB2	1:A:301:PRO:HD3	2.02	0.40
1:B:293:ASP:OD2	1:B:316:TYR:OH	2.29	0.40
1:A:327:ARG:CZ	1:A:364:ARG:HD2	2.51	0.40
1:B:347:HIS:CG	1:B:348:GLY:N	2.88	0.40
1:A:155:TYR:CE1	1:D:461:GLY:HA3	2.57	0.40
1:A:238:THR:HG21	1:A:287:VAL:HG21	2.03	0.40
1:A:227:GLU:O	1:A:350:LYS:HE2	2.22	0.40
1:C:166:PRO:O	1:C:168:HIS:HD2	2.04	0.40
1:C:194:ILE:HA	1:C:202:THR:O	2.22	0.40
1:A:461:GLY:HA3	1:B:155:TYR:CE1	2.57	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	386/393 (98%)	358 (93%)	25 (6%)	3 (1%)	24	22
1	B	386/393 (98%)	362 (94%)	21 (5%)	3 (1%)	24	22
1	C	386/393 (98%)	361 (94%)	23 (6%)	2 (0%)	34	35
1	D	386/393 (98%)	362 (94%)	22 (6%)	2 (0%)	34	35
All	All	1544/1572 (98%)	1443 (94%)	91 (6%)	10 (1%)	30	29

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	468	PRO
1	A	468	PRO

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Mol	Chain	Res	Type
1	C	234	ASN
1	A	329	ASN
1	B	468	PRO
1	D	319	SER
1	D	468	PRO
1	B	319	SER
1	A	319	SER
1	B	336	HIS

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	337/341 (99%)	320 (95%)	17 (5%)	30	35
1	B	337/341 (99%)	326 (97%)	11 (3%)	45	56
1	C	337/341 (99%)	321 (95%)	16 (5%)	32	39
1	D	337/341 (99%)	328 (97%)	9 (3%)	52	64
All	All	1348/1364 (99%)	1295 (96%)	53 (4%)	39	48

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	ILE
1	A	105	SER
1	A	128	LYS
1	A	216	VAL
1	A	247	SER
1	A	249	LYS
1	A	267	THR
1	A	276	GLU
1	A	359	ASP
1	A	368	GLU
1	A	372	LEU
1	A	385	ASN
1	A	399	ASP

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Mol	Chain	Res	Type
1	A	416	SER
1	A	435	GLU
1	A	467	MET
1	A	469	ILE
1	B	94	ILE
1	B	216	VAL
1	B	267	THR
1	B	307	ILE
1	B	368	GLU
1	B	385	ASN
1	B	415	LYS
1	B	431	LYS
1	B	435	GLU
1	B	463	ASP
1	B	468	PRO
1	C	105	SER
1	C	215	VAL
1	C	216	VAL
1	C	267	THR
1	C	276	GLU
1	C	307	ILE
1	C	330	ASP
1	C	372	LEU
1	C	385	ASN
1	C	388	SER
1	C	400	ARG
1	C	431	LYS
1	C	434	THR
1	C	463	ASP
1	C	467	MET
1	C	468	PRO
1	D	105	SER
1	D	128	LYS
1	D	216	VAL
1	D	267	THR
1	D	307	ILE
1	D	334	SER
1	D	370	SER
1	D	385	ASN
1	D	416	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	136	GLN
1	A	385	ASN
1	A	391	GLN
1	A	393	ASN
1	B	136	GLN
1	B	385	ASN
1	B	393	ASN
1	B	402	ASN
1	C	136	GLN
1	C	273	GLN
1	C	347	HIS
1	C	385	ASN
1	C	393	ASN
1	D	136	GLN
1	D	168	HIS
1	D	358	ASN
1	D	385	ASN
1	D	393	ASN
1	D	402	ASN
1	D	465	ASN

5.3.3 RNA [i](#)

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](#)

28 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	NAG	A	502	1,3	14,14,15	0.92	1 (7%)	15,19,21	2.26	2 (13%)
3	NAG	A	503	3	14,14,15	0.63	0	15,19,21	1.58	3 (20%)
3	BMA	A	504	3	11,11,12	0.99	1 (9%)	14,15,17	1.82	4 (28%)
3	BMA	A	505	3	11,11,12	1.19	2 (18%)	14,15,17	5.09	8 (57%)
3	BMA	A	506	3	11,11,12	0.47	0	14,15,17	3.60	6 (42%)
4	NAG	A	507	1,4	14,14,15	0.72	0	15,19,21	1.35	3 (20%)
4	NAG	A	508	4	14,14,15	0.58	0	15,19,21	1.21	2 (13%)
3	NAG	B	502	1,3	14,14,15	0.93	0	15,19,21	1.49	2 (13%)
3	NAG	B	503	3	14,14,15	0.74	0	15,19,21	1.88	4 (26%)
3	BMA	B	504	3	11,11,12	1.11	1 (9%)	14,15,17	1.28	2 (14%)
3	BMA	B	505	3	11,11,12	1.15	2 (18%)	14,15,17	4.24	5 (35%)
3	BMA	B	506	3	11,11,12	0.70	0	14,15,17	3.70	8 (57%)
3	NAG	C	502	1,3	14,14,15	0.92	1 (7%)	15,19,21	1.89	2 (13%)
3	NAG	C	503	3	14,14,15	0.78	0	15,19,21	1.88	4 (26%)
3	BMA	C	504	3	11,11,12	0.72	0	14,15,17	2.02	3 (21%)
3	BMA	C	505	3	11,11,12	1.19	2 (18%)	14,15,17	4.50	6 (42%)
3	BMA	C	506	3	11,11,12	0.87	0	14,15,17	3.14	7 (50%)
7	NAG	C	507	1,7	14,14,15	0.57	0	15,19,21	1.45	2 (13%)
7	FUC	C	508	7	10,10,11	0.76	0	14,14,16	1.52	3 (21%)
3	NAG	D	502	1,3	14,14,15	0.97	0	15,19,21	2.16	5 (33%)
3	NAG	D	503	3	14,14,15	0.69	0	15,19,21	1.75	5 (33%)
3	BMA	D	504	3	11,11,12	0.73	0	14,15,17	1.90	4 (28%)
3	BMA	D	505	3	11,11,12	1.05	1 (9%)	14,15,17	4.42	7 (50%)
3	BMA	D	506	3	11,11,12	0.58	0	14,15,17	3.49	8 (57%)
8	NAG	D	508	1,8	14,14,15	0.56	0	15,19,21	1.52	4 (26%)
8	NAG	D	509	8	14,14,15	0.54	0	15,19,21	1.17	1 (6%)
8	FUC	D	510	8	10,10,11	1.02	0	14,14,16	2.10	4 (28%)
8	FUC	D	511	8	10,10,11	0.65	0	14,14,16	2.26	6 (42%)

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	NAG	A	502	1,3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	503	3	-	0/6/23/26	0/1/1/1
3	BMA	A	504	3	-	0/2/19/22	0/1/1/1
3	BMA	A	505	3	-	0/2/19/22	0/1/1/1
3	BMA	A	506	3	-	0/2/19/22	0/1/1/1
4	NAG	A	507	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	508	4	-	0/6/23/26	0/1/1/1
3	NAG	B	502	1,3	-	0/6/23/26	0/1/1/1
3	NAG	B	503	3	-	0/6/23/26	0/1/1/1
3	BMA	B	504	3	-	0/2/19/22	0/1/1/1
3	BMA	B	505	3	-	0/2/19/22	0/1/1/1
3	BMA	B	506	3	-	0/2/19/22	0/1/1/1
3	NAG	C	502	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	503	3	-	0/6/23/26	0/1/1/1
3	BMA	C	504	3	-	0/2/19/22	0/1/1/1
3	BMA	C	505	3	-	0/2/19/22	0/1/1/1
3	BMA	C	506	3	-	0/2/19/22	0/1/1/1
7	NAG	C	507	1,7	-	0/6/23/26	0/1/1/1
7	FUC	C	508	7	-	0/0/17/20	0/1/1/1
3	NAG	D	502	1,3	-	0/6/23/26	0/1/1/1
3	NAG	D	503	3	-	0/6/23/26	0/1/1/1
3	BMA	D	504	3	-	0/2/19/22	0/1/1/1
3	BMA	D	505	3	-	0/2/19/22	0/1/1/1
3	BMA	D	506	3	-	0/2/19/22	0/1/1/1
8	NAG	D	508	1,8	-	0/6/23/26	0/1/1/1
8	NAG	D	509	8	-	0/6/23/26	0/1/1/1
8	FUC	D	510	8	-	0/0/17/20	0/1/1/1
8	FUC	D	511	8	-	0/0/17/20	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	505	BMA	O5-C1	-2.33	1.39	1.43
3	B	505	BMA	O5-C1	-2.20	1.40	1.43
3	C	502	NAG	O5-C1	-2.14	1.40	1.43
3	A	505	BMA	O5-C1	-2.12	1.40	1.43
3	B	504	BMA	O5-C1	-2.12	1.40	1.43
3	A	502	NAG	O3-C3	-2.10	1.37	1.43
3	D	505	BMA	C2-C3	2.07	1.55	1.52
3	B	505	BMA	C2-C3	2.14	1.55	1.52
3	C	505	BMA	C2-C3	2.19	1.55	1.52
3	A	505	BMA	C4-C5	2.35	1.58	1.53
3	A	504	BMA	C2-C3	2.54	1.56	1.52

All (120) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	505	BMA	C1-C2-C3	-13.94	93.05	109.54
3	C	505	BMA	C1-C2-C3	-11.90	95.47	109.54
3	D	505	BMA	C1-C2-C3	-11.84	95.53	109.54
3	B	505	BMA	C1-C2-C3	-11.59	95.84	109.54
3	B	506	BMA	C1-O5-C5	-11.10	98.16	112.25
3	A	506	BMA	C1-O5-C5	-9.99	99.57	112.25
3	C	505	BMA	C1-O5-C5	-8.83	101.04	112.25
3	A	505	BMA	C1-O5-C5	-8.65	101.27	112.25
3	D	505	BMA	C1-O5-C5	-8.29	101.73	112.25
3	B	505	BMA	C1-O5-C5	-7.97	102.13	112.25
3	C	506	BMA	C1-O5-C5	-7.22	103.08	112.25
3	C	506	BMA	C1-C2-C3	-5.85	102.62	109.54
8	D	510	FUC	O5-C1-C2	-5.07	102.63	110.86
3	C	504	BMA	C1-O5-C5	-4.80	106.16	112.25
3	D	504	BMA	O3-C3-C4	-4.58	100.02	110.34
3	A	502	NAG	O3-C3-C4	-4.44	100.34	110.34
3	D	502	NAG	C2-N2-C7	-4.30	117.52	123.04
3	B	503	NAG	O4-C4-C3	-4.08	101.14	110.34
3	C	504	BMA	O3-C3-C4	-3.90	101.56	110.34
3	B	506	BMA	O4-C4-C3	-3.85	101.67	110.34
3	B	503	NAG	C3-C2-N2	-3.73	101.62	110.56
3	B	506	BMA	C1-C2-C3	-3.70	105.16	109.54
3	D	506	BMA	O4-C4-C3	-3.68	102.04	110.34
3	C	503	NAG	C3-C2-N2	-3.67	101.76	110.56
3	D	503	NAG	O4-C4-C3	-3.29	102.93	110.34
3	A	504	BMA	O3-C3-C4	-3.29	102.94	110.34
3	A	506	BMA	O4-C4-C3	-3.27	102.98	110.34
3	A	504	BMA	C2-C3-C4	-3.08	105.81	111.04
7	C	507	NAG	O6-C6-C5	-2.95	101.60	111.33
3	D	502	NAG	O7-C7-C8	-2.95	116.66	122.06
3	A	504	BMA	O2-C2-C1	-2.92	103.36	109.21
3	D	503	NAG	O7-C7-C8	-2.88	116.77	122.06
4	A	508	NAG	C3-C4-C5	-2.88	105.18	110.20
3	A	503	NAG	C3-C2-N2	-2.86	103.70	110.56
3	D	506	BMA	O3-C3-C4	-2.83	103.96	110.34
7	C	507	NAG	C2-N2-C7	-2.80	119.44	123.04
8	D	511	FUC	O3-C3-C2	-2.80	104.94	110.00
3	C	504	BMA	O6-C6-C5	-2.78	102.14	111.33
3	D	503	NAG	C3-C2-N2	-2.64	104.22	110.56
8	D	508	NAG	O6-C6-C5	-2.64	102.59	111.33
3	A	503	NAG	O6-C6-C5	-2.56	102.89	111.33
3	D	504	BMA	C1-O5-C5	-2.53	109.04	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	NAG	O3-C3-C4	-2.50	104.71	110.34
3	C	503	NAG	O4-C4-C3	-2.49	104.73	110.34
3	C	502	NAG	C2-N2-C7	-2.49	119.84	123.04
4	A	507	NAG	C3-C2-N2	-2.40	104.81	110.56
3	A	503	NAG	O7-C7-C8	-2.34	117.77	122.06
3	B	503	NAG	C3-C4-C5	-2.14	106.46	110.20
4	A	507	NAG	O6-C6-C5	-2.14	104.27	111.33
3	D	506	BMA	O6-C6-C5	-2.11	104.36	111.33
3	B	504	BMA	O3-C3-C4	-2.08	105.66	110.34
8	D	511	FUC	O2-C2-C3	-2.03	106.03	110.12
3	B	506	BMA	O3-C3-C4	-2.02	105.78	110.34
4	A	507	NAG	C2-N2-C7	-2.02	120.44	123.04
3	D	505	BMA	O5-C5-C6	2.01	111.70	107.35
3	D	503	NAG	C6-C5-C4	2.02	118.00	113.02
3	C	506	BMA	C6-C5-C4	2.04	118.05	113.02
3	B	506	BMA	O5-C5-C6	2.05	111.79	107.35
3	B	503	NAG	C1-O5-C5	2.07	114.87	112.25
8	D	511	FUC	O5-C5-C6	2.08	109.57	106.13
3	D	504	BMA	O3-C3-C2	2.10	113.79	110.00
3	C	503	NAG	C8-C7-N2	2.11	120.14	116.11
3	A	505	BMA	O2-C2-C3	2.15	114.45	110.12
3	D	502	NAG	C8-C7-N2	2.17	120.27	116.11
3	B	505	BMA	O3-C3-C2	2.21	113.99	110.00
3	D	506	BMA	O2-C2-C1	2.28	113.78	109.21
7	C	508	FUC	O5-C5-C4	2.29	113.49	109.53
8	D	511	FUC	C3-C4-C5	2.34	113.66	109.72
3	C	506	BMA	O2-C2-C3	2.35	114.84	110.12
4	A	508	NAG	O5-C5-C6	2.36	112.45	107.35
8	D	508	NAG	O5-C5-C6	2.36	112.45	107.35
3	B	506	BMA	O4-C4-C5	2.36	115.49	109.24
8	D	510	FUC	O5-C5-C6	2.36	110.03	106.13
3	D	505	BMA	O2-C2-C3	2.37	114.88	110.12
7	C	508	FUC	C1-O5-C5	2.39	116.08	112.38
3	D	504	BMA	O5-C5-C6	2.43	112.61	107.35
3	C	506	BMA	C2-C3-C4	2.44	115.18	111.04
3	D	502	NAG	O5-C5-C6	2.47	112.70	107.35
3	B	504	BMA	C1-O5-C5	2.56	115.50	112.25
3	B	506	BMA	O5-C1-C2	2.59	115.06	110.86
3	C	505	BMA	O2-C2-C3	2.62	115.39	110.12
8	D	508	NAG	C4-C3-C2	2.63	115.32	111.23
3	A	505	BMA	O4-C4-C5	2.67	116.32	109.24
3	D	503	NAG	C8-C7-N2	2.67	121.22	116.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	504	BMA	O3-C3-C2	2.70	114.87	110.00
3	C	505	BMA	O3-C3-C2	2.83	115.11	110.00
8	D	510	FUC	C3-C4-C5	2.84	114.51	109.72
3	D	505	BMA	O3-C3-C2	2.92	115.28	110.00
3	A	506	BMA	C2-C3-C4	2.96	116.08	111.04
8	D	509	NAG	C1-O5-C5	2.99	116.04	112.25
3	A	505	BMA	O3-C3-C4	3.02	117.13	110.34
8	D	508	NAG	C1-O5-C5	3.04	116.10	112.25
3	A	506	BMA	O5-C1-C2	3.07	115.84	110.86
3	D	506	BMA	C3-C4-C5	3.16	115.70	110.20
3	B	506	BMA	C2-C3-C4	3.21	116.50	111.04
8	D	511	FUC	C2-C3-C4	3.29	116.64	111.04
3	D	505	BMA	O5-C1-C2	3.40	116.37	110.86
7	C	508	FUC	O5-C5-C6	3.43	111.81	106.13
3	C	505	BMA	O5-C1-C2	3.49	116.53	110.86
3	D	506	BMA	O5-C1-C2	3.56	116.62	110.86
3	B	505	BMA	O5-C1-C2	3.65	116.78	110.86
8	D	510	FUC	O5-C5-C4	3.67	115.88	109.53
3	B	502	NAG	C1-O5-C5	3.77	117.04	112.25
3	C	506	BMA	O2-C2-C1	3.83	116.89	109.21
3	D	506	BMA	C2-C3-C4	3.87	117.61	111.04
3	C	506	BMA	O5-C1-C2	3.97	117.30	110.86
3	C	503	NAG	C1-O5-C5	4.37	117.79	112.25
3	A	505	BMA	C6-C5-C4	4.42	123.91	113.02
3	B	505	BMA	O2-C2-C1	4.64	118.52	109.21
3	A	506	BMA	O5-C5-C6	4.74	117.61	107.35
3	A	505	BMA	O5-C1-C2	4.80	118.65	110.86
3	D	502	NAG	C1-O5-C5	4.92	118.49	112.25
3	A	506	BMA	C1-C2-C3	5.04	115.50	109.54
3	A	505	BMA	O2-C2-C1	5.10	119.44	109.21
3	C	505	BMA	O2-C2-C1	5.13	119.50	109.21
3	D	505	BMA	O2-C2-C1	5.27	119.77	109.21
8	D	511	FUC	C1-C2-C3	5.68	116.26	109.54
3	C	502	NAG	C1-O5-C5	6.19	120.11	112.25
3	A	502	NAG	C1-O5-C5	7.12	121.29	112.25
3	D	506	BMA	C1-O5-C5	9.76	124.64	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	508	FUC	1	0
3	D	506	BMA	1	0
8	D	510	FUC	1	0

5.6 Ligand geometry

Of 17 ligands modelled in this entry, 5 are monoatomic - leaving 12 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	G39	A	501	-	16,20,20	1.36	3 (18%)	12,27,27	1.10	0
5	NAG	A	509	1	14,14,15	0.81	1 (7%)	15,19,21	1.73	1 (6%)
5	NAG	A	510	1	14,14,15	0.47	0	15,19,21	1.64	1 (6%)
2	G39	B	501	-	16,20,20	0.97	1 (6%)	12,27,27	1.59	2 (16%)
5	NAG	B	507	1	14,14,15	0.53	0	15,19,21	1.61	2 (13%)
5	NAG	B	508	1	14,14,15	0.58	0	15,19,21	1.96	3 (20%)
5	NAG	B	509	1	14,14,15	0.49	0	15,19,21	2.51	3 (20%)
2	G39	C	501	-	16,20,20	1.31	3 (18%)	12,27,27	0.81	0
5	NAG	C	509	1	14,14,15	0.49	0	15,19,21	1.84	2 (13%)
2	G39	D	501	-	16,20,20	1.48	3 (18%)	12,27,27	1.33	2 (16%)
5	NAG	D	507	1	14,14,15	0.59	0	15,19,21	2.27	1 (6%)
5	NAG	D	512	1	14,14,15	0.54	0	15,19,21	1.79	4 (26%)

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	G39	A	501	-	-	0/12/32/32	0/1/1/1
5	NAG	A	509	1	-	0/6/23/26	0/1/1/1
5	NAG	A	510	1	-	0/6/23/26	0/1/1/1
2	G39	B	501	-	-	0/12/32/32	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	B	507	1	-	0/6/23/26	0/1/1/1
5	NAG	B	508	1	-	0/6/23/26	0/1/1/1
5	NAG	B	509	1	-	0/6/23/26	0/1/1/1
2	G39	C	501	-	-	0/12/32/32	0/1/1/1
5	NAG	C	509	1	-	0/6/23/26	0/1/1/1
2	G39	D	501	-	-	0/12/32/32	0/1/1/1
5	NAG	D	507	1	-	0/6/23/26	0/1/1/1
5	NAG	D	512	1	-	0/6/23/26	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	G39	C1-C2	-3.81	1.44	1.51
2	D	501	G39	C1-C2	-2.94	1.46	1.51
2	A	501	G39	C11-C10	-2.62	1.45	1.50
2	D	501	G39	C11-C10	-2.58	1.45	1.50
2	C	501	G39	C5-N5	-2.29	1.42	1.45
2	B	501	G39	C3-C2	2.07	1.53	1.50
2	A	501	G39	C3-C4	2.09	1.58	1.54
2	C	501	G39	C3-C2	2.38	1.54	1.50
5	A	509	NAG	C1-C2	2.40	1.55	1.52
2	A	501	G39	C3-C2	2.79	1.55	1.50
2	D	501	G39	C3-C2	3.48	1.56	1.50

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	G39	C6-C7-C2	-3.53	116.81	122.39
5	D	512	NAG	C4-C3-C2	-3.42	105.91	111.23
5	B	509	NAG	C4-C3-C2	-2.79	106.90	111.23
5	B	508	NAG	C2-N2-C7	-2.55	119.76	123.04
5	D	512	NAG	C2-N2-C7	-2.52	119.80	123.04
5	B	508	NAG	C3-C4-C5	-2.21	106.34	110.20
2	D	501	G39	C5-N5-C10	-2.20	117.45	123.10
5	C	509	NAG	C4-C3-C2	2.09	114.48	111.23
5	D	512	NAG	O4-C4-C5	2.10	114.80	109.24
2	B	501	G39	C4-C3-C2	2.27	112.54	109.65
2	D	501	G39	O10-C10-C11	2.58	126.79	122.06
5	B	509	NAG	C2-N2-C7	2.61	126.40	123.04
5	B	507	NAG	C2-N2-C7	3.40	127.41	123.04
5	D	512	NAG	C1-O5-C5	3.48	116.67	112.25
5	B	507	NAG	C1-O5-C5	3.85	117.13	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	510	NAG	C1-O5-C5	5.53	119.26	112.25
5	A	509	NAG	C1-O5-C5	5.56	119.30	112.25
5	C	509	NAG	C1-O5-C5	5.66	119.44	112.25
5	B	508	NAG	C1-O5-C5	6.00	119.86	112.25
5	B	509	NAG	C1-O5-C5	8.09	122.51	112.25
5	D	507	NAG	C1-O5-C5	8.30	122.78	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	G39	3	0
5	A	510	NAG	1	0
2	B	501	G39	2	0
5	C	509	NAG	1	0
2	D	501	G39	1	0
5	D	512	NAG	1	0

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, 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	388/393 (98%)	0.24	12 (3%)	52 51	19, 31, 49, 77	0
1	B	388/393 (98%)	0.27	19 (4%)	33 33	18, 31, 50, 74	0
1	C	388/393 (98%)	0.26	16 (4%)	41 39	17, 31, 49, 78	0
1	D	388/393 (98%)	0.15	16 (4%)	41 39	19, 31, 48, 75	0
All	All	1552/1572 (98%)	0.23	63 (4%)	41 39	17, 31, 49, 78	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	469	ILE	7.2
1	A	469	ILE	5.9
1	D	468	PRO	5.9
1	C	469	ILE	5.5
1	D	469	ILE	5.4
1	C	330	ASP	5.0
1	A	468	PRO	4.5
1	C	468	PRO	4.3
1	B	82	ALA	4.1
1	B	468	PRO	4.0
1	A	169	LEU	3.3
1	D	332	SER	3.3
1	A	390	LEU	3.3
1	D	331	SER	3.2
1	A	331	SER	3.0
1	D	159	LEU	2.9
1	B	83	GLU	2.9
1	B	165	VAL	2.9
1	B	169	LEU	2.8
1	C	169	LEU	2.8
1	A	171	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	332	SER	2.8
1	B	331	SER	2.7
1	B	390	LEU	2.7
1	C	332	SER	2.7
1	B	159	LEU	2.6
1	A	321	LEU	2.6
1	B	174	VAL	2.6
1	B	168	HIS	2.6
1	B	171	THR	2.5
1	A	114	ILE	2.5
1	A	330	ASP	2.5
1	D	174	VAL	2.5
1	D	334	SER	2.5
1	B	166	PRO	2.5
1	C	114	ILE	2.4
1	C	368	GLU	2.4
1	C	174	VAL	2.4
1	D	169	LEU	2.4
1	D	330	ASP	2.4
1	B	431	LYS	2.3
1	B	167	PHE	2.3
1	C	168	HIS	2.3
1	D	114	ILE	2.3
1	B	170	GLY	2.2
1	C	247	SER	2.2
1	C	344	GLU	2.2
1	D	171	THR	2.2
1	D	167	PHE	2.2
1	C	331	SER	2.2
1	C	392	ILE	2.2
1	A	163	LEU	2.2
1	D	168	HIS	2.2
1	C	165	VAL	2.2
1	D	165	VAL	2.2
1	B	334	SER	2.1
1	D	328	LYS	2.1
1	D	161	ASN	2.1
1	A	168	HIS	2.1
1	B	321	LEU	2.0
1	C	159	LEU	2.0
1	A	328	LYS	2.0
1	C	86	ASN	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
3	NAG	D	502	14/15	0.95	0.15	2.99	29,32,44,44	0
3	NAG	B	502	14/15	0.96	0.12	1.98	30,34,47,49	0
3	NAG	A	502	14/15	0.94	0.13	1.65	28,36,45,47	0
3	NAG	C	502	14/15	0.96	0.11	0.24	31,35,46,47	0
3	BMA	D	506	11/12	0.89	0.13	-	55,62,66,67	0
8	NAG	D	509	14/15	0.87	0.28	-	59,70,75,75	0
3	BMA	A	504	11/12	0.93	0.16	-	38,43,48,54	0
3	BMA	B	505	11/12	0.91	0.20	-	39,48,55,61	0
3	NAG	C	503	14/15	0.97	0.12	-	31,34,40,41	0
3	BMA	B	504	11/12	0.92	0.18	-	38,40,48,61	0
3	BMA	A	506	11/12	0.84	0.15	-	54,58,65,65	0
3	BMA	D	505	11/12	0.83	0.20	-	46,51,55,55	0
4	NAG	A	507	14/15	0.83	0.15	-	54,64,74,80	0
3	BMA	C	505	11/12	0.90	0.17	-	45,49,53,53	0
3	BMA	A	505	11/12	0.82	0.20	-	45,46,55,55	0
8	NAG	D	508	14/15	0.93	0.14	-	46,51,59,60	0
4	NAG	A	508	14/15	0.81	0.29	-	74,93,104,109	0
3	NAG	D	503	14/15	0.95	0.14	-	31,35,41,44	0
3	BMA	B	506	11/12	0.82	0.20	-	46,50,55,62	0
3	NAG	B	503	14/15	0.97	0.11	-	31,36,38,40	0
3	NAG	A	503	14/15	0.94	0.12	-	30,36,37,38	0
7	NAG	C	507	14/15	0.87	0.15	-	49,53,63,69	0
3	BMA	C	506	11/12	0.91	0.15	-	42,49,56,58	0
3	BMA	D	504	11/12	0.90	0.12	-	38,45,51,54	0
8	FUC	D	510	10/11	0.81	0.26	-	66,72,76,82	0
7	FUC	C	508	10/11	0.89	0.32	-	74,77,82,83	0
3	BMA	C	504	11/12	0.89	0.15	-	40,42,47,51	0
8	FUC	D	511	10/11	0.91	0.31	-	60,66,71,74	0

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	G39	D	501	20/20	0.94	0.16	1.54	23,29,33,47	0
2	G39	B	501	20/20	0.94	0.15	1.38	26,29,32,40	0
2	G39	A	501	20/20	0.92	0.15	1.27	26,30,35,49	0
2	G39	C	501	20/20	0.95	0.13	0.10	25,29,37,39	0
5	NAG	A	510	14/15	0.88	0.17	-0.56	55,60,66,67	0
6	CA	D	513	1/1	0.98	0.04	-1.96	33,33,33,33	0
6	CA	A	511	1/1	0.98	0.04	-2.26	34,34,34,34	0
6	CA	C	510	1/1	0.97	0.04	-2.34	38,38,38,38	0
6	CA	B	510	1/1	0.99	0.03	-2.57	35,35,35,35	0
6	CA	A	512	1/1	0.96	0.05	-	61,61,61,61	0
5	NAG	D	512	14/15	0.87	0.27	-	54,59,69,73	0
5	NAG	B	508	14/15	0.89	0.15	-	58,64,68,74	0
5	NAG	A	509	14/15	0.89	0.35	-	49,63,71,77	0
5	NAG	B	507	14/15	0.83	0.27	-	61,68,75,77	0
5	NAG	B	509	14/15	0.85	0.24	-	57,66,72,72	0
5	NAG	D	507	14/15	0.78	0.32	-	56,70,75,76	0
5	NAG	C	509	14/15	0.89	0.26	-	62,69,71,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.