



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:45 PM GMT

PDB ID : 1V0Z
Title : STRUCTURE OF NEURAMINIDASE FROM ENGLISH DUCK SUBTYPE N6
Authors : Rudino-Pinera, E.; Tunnah, P.; Crennell, S.J.; Webster, R.G.; Laver, W.G.; Garman, E.F.
Deposited on : 2004-04-02
Resolution : 1.84 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

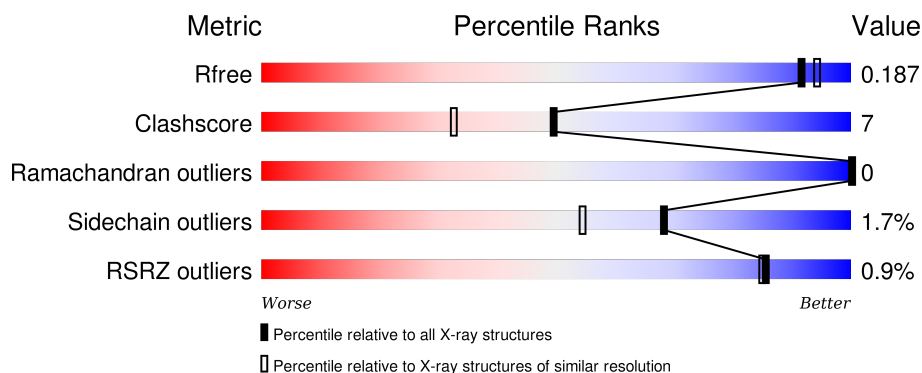
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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	2634 (1.86-1.82)
Clashscore	102246	2862 (1.86-1.82)
Ramachandran outliers	100387	2831 (1.86-1.82)
Sidechain outliers	100360	2832 (1.86-1.82)
RSRZ outliers	91569	2639 (1.86-1.82)

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	389	<div> <div>%</div> <div> <div></div> <div>91%</div> <div>8%</div> <div>.</div> </div> </div>
1	B	389	<div> <div></div> <div>92%</div> <div>7%</div> <div>.</div> </div>
1	C	389	<div> <div>2%</div> <div></div> <div>89%</div> <div>10%</div> </div>
1	D	389	<div> <div>%</div> <div></div> <div>89%</div> <div>10%</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
10	PEG	C	1478	-	-	X	X
2	CA	C	1477	-	-	-	X
2	CA	D	1477	-	-	-	X
3	GOL	A	1478	-	-	-	X
3	GOL	B	1478	-	-	-	X
3	GOL	C	1479	-	-	-	X
3	GOL	D	1478	-	-	-	X
4	NAG	C	1483	-	-	-	X
5	NAG	B	1482	-	-	X	X
5	NAG	D	1481	-	-	X	X
6	MAN	A	1483	-	-	-	X
6	NAG	A	1495	-	-	-	X
7	MAN	A	1485	-	-	-	X
7	MAN	A	1494	-	-	-	X
7	MAN	B	1485	-	-	-	X
7	MAN	C	1485	-	-	-	X
8	NAG	A	1488	-	-	-	X
8	MAN	A	1493	-	-	-	X
9	NAG	B	1487	-	-	-	X

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 14347 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	389	Total	C	N	O	S	0	0	0
			3009	1875	535	572	27			
1	B	389	Total	C	N	O	S	0	0	0
			3009	1875	535	572	27			
1	C	389	Total	C	N	O	S	0	0	0
			3009	1875	535	572	27			
1	D	389	Total	C	N	O	S	0	0	0
			3009	1875	535	572	27			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- 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		
4	C	2	Total	C	N	O	0	0
			28	16	2	10		

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

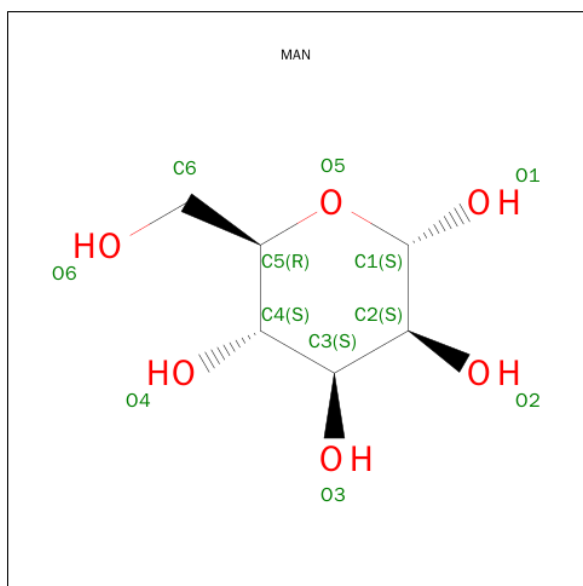


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		
5	C	1	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 a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	4	Total	C	N	O	0	0
			47	26	1	20		

- Molecule 7 is SUGAR (ALPHA-D-MANNOSE) (three-letter code: MAN) (formula: $C_6H_{12}O_6$).



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

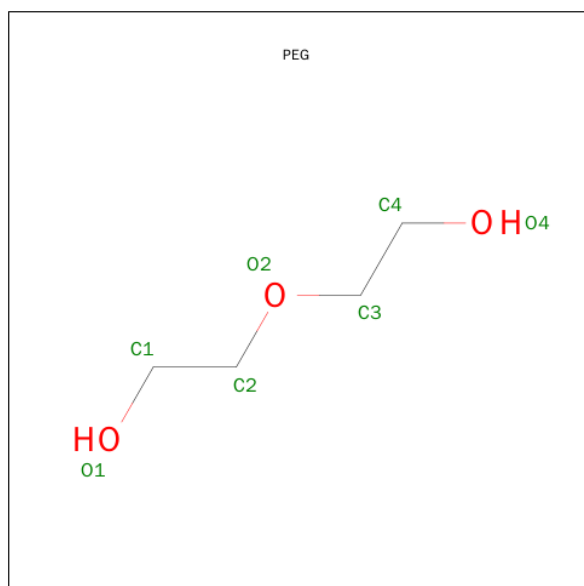
- Molecule 8 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 9 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	3	Total	C	N	O	
			36	20	1	15	
						0	0

- Molecule 10 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	C	1	Total	C	O		
			7	4	3		
						0	0

- Molecule 11 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	C	2	Total	C	O		
			22	12	10		
						0	0

- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	453	Total	O		
			453	453	0	0
12	B	434	Total	O		
			434	434	0	0
12	C	450	Total	O		
			450	450	0	0

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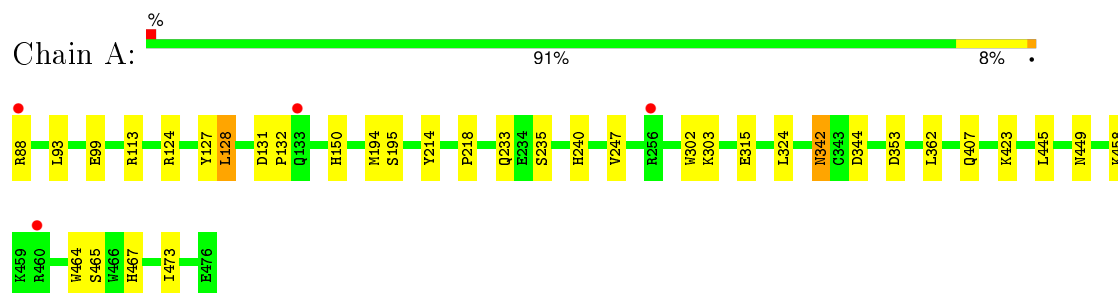
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	D	461	Total 461	O 461	0	0

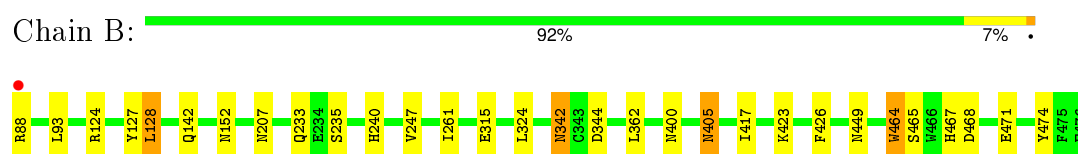
3 Residue-property plots [i](#)

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.

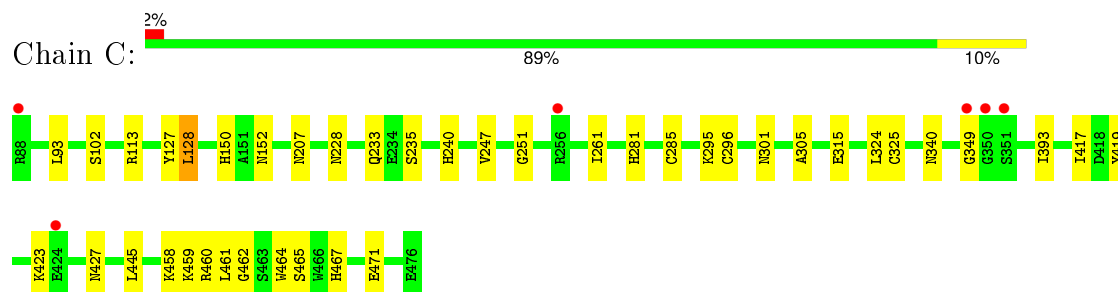
• Molecule 1: NEURAMINIDASE



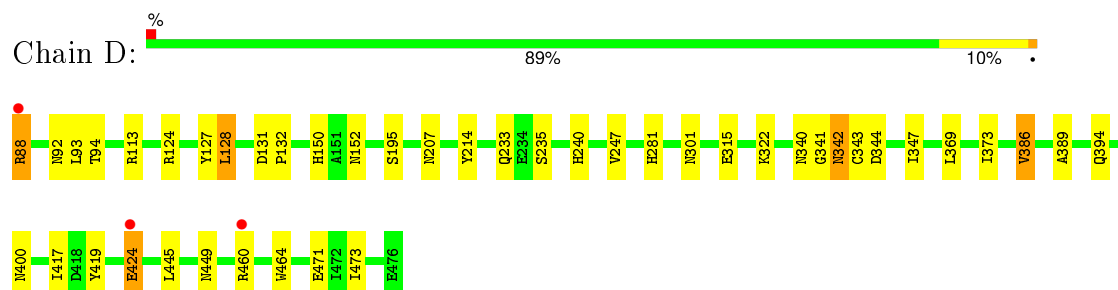
• Molecule 1: NEURAMINIDASE



• 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	106.48Å 73.75Å 106.81Å 90.00° 90.37° 90.00°	Depositor
Resolution (Å)	19.96 – 1.84 19.98 – 1.84	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.96-1.84) 93.2 (19.98-1.84)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 1.84Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.149 , 0.190 0.150 , 0.187	Depositor DCC
R_{free} test set	6696 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	12.7	Xtriage
Anisotropy	1.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 61.4	EDS
Estimated twinning fraction	0.006 for l,k,-h 0.016 for h,-k,-l 0.012 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 133791 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14347	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.12% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, BMA, NAG, CA, PEG, 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	A	0.65	0/3084	0.67	1/4185 (0.0%)
1	B	0.65	1/3084 (0.0%)	0.67	1/4185 (0.0%)
1	C	0.66	0/3084	0.66	1/4185 (0.0%)
1	D	0.65	0/3084	0.69	1/4185 (0.0%)
All	All	0.65	1/12336 (0.0%)	0.67	4/16740 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	464	TRP	CB-CG	-5.11	1.41	1.50

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	128	LEU	CA-CB-CG	7.07	131.57	115.30
1	D	128	LEU	CA-CB-CG	6.70	130.71	115.30
1	B	128	LEU	CA-CB-CG	6.39	129.99	115.30
1	A	128	LEU	CA-CB-CG	6.12	129.38	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3009	0	2885	26	0
1	B	3009	0	2888	30	0
1	C	3009	0	2888	51	0
1	D	3009	0	2888	48	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	6	0	8	2	0
3	B	6	0	8	1	0
3	C	6	0	8	1	0
3	D	6	0	8	0	0
4	A	28	0	25	0	0
4	C	28	0	25	6	0
5	A	28	0	26	0	0
5	B	56	0	52	11	0
5	C	42	0	39	5	0
5	D	42	0	39	13	0
6	A	47	0	40	4	0
7	A	33	0	30	4	0
7	B	22	0	20	0	0
7	C	11	0	10	1	0
7	D	11	0	10	2	0
8	A	72	0	61	6	0
9	B	36	0	31	5	0
10	C	7	0	10	16	0
11	C	22	0	19	5	0
12	A	453	0	0	5	0
12	B	434	0	0	1	0
12	C	450	0	0	6	0
12	D	461	0	0	1	0
All	All	14347	0	12018	163	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 7.

All (163) 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:D:207:ASN:HD21	5:D:1481:NAG:C1	1.03	1.63
1:D:152:ASN:HD21	5:D:1480:NAG:C1	1.01	1.55
1:B:152:ASN:HD21	5:B:1480:NAG:C1	0.97	1.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:ASN:HD21	5:C:1481:NAG:C1	0.92	1.53
1:C:207:ASN:HD21	4:C:1483:NAG:C1	0.94	1.52
1:C:460:ARG:NH1	12:C:2412:HOH:O	1.60	1.27
1:D:88:ARG:HB2	1:D:88:ARG:HH11	0.98	1.08
1:D:88:ARG:HH11	1:D:88:ARG:CB	1.73	1.02
1:C:423:LYS:NZ	10:C:1478:PEG:H32	1.76	1.00
1:C:423:LYS:NZ	10:C:1478:PEG:C3	2.27	0.98
1:B:93:LEU:H	1:B:240:HIS:HD2	1.12	0.98
1:D:88:ARG:HB2	1:D:88:ARG:NH1	1.78	0.97
1:C:93:LEU:H	1:C:240:HIS:HD2	1.13	0.95
9:B:1487:NAG:C1	5:D:1481:NAG:O4	2.15	0.93
6:A:1495:NAG:C1	5:B:1482:NAG:HO4	1.78	0.92
1:C:423:LYS:NZ	10:C:1478:PEG:H42	1.83	0.92
1:C:150:HIS:HE1	1:D:471:GLU:H	1.13	0.92
1:B:471:GLU:H	1:D:150:HIS:HE1	1.15	0.90
1:D:93:LEU:H	1:D:240:HIS:HD2	1.17	0.90
1:A:150:HIS:HE1	1:C:471:GLU:H	1.17	0.88
1:B:423:LYS:HE2	12:D:2188:HOH:O	1.75	0.86
4:C:1484:NAG:O4	11:C:1486:BMA:C1	2.24	0.84
1:B:207:ASN:HD21	5:B:1482:NAG:C1	1.90	0.84
1:B:93:LEU:H	1:B:240:HIS:CD2	1.98	0.82
1:B:468:ASP:O	12:B:2402:HOH:O	1.97	0.82
9:B:1487:NAG:C1	5:D:1481:NAG:C4	2.59	0.81
1:C:423:LYS:NZ	10:C:1478:PEG:C4	2.44	0.80
1:A:93:LEU:H	1:A:240:HIS:HD2	1.27	0.80
1:C:423:LYS:HZ1	10:C:1478:PEG:C4	1.96	0.78
1:C:423:LYS:HZ1	10:C:1478:PEG:C3	1.95	0.78
1:D:386:VAL:HG22	1:D:389:ALA:HB2	1.66	0.77
1:C:423:LYS:HZ1	10:C:1478:PEG:H42	1.49	0.76
1:C:233:GLN:HE21	1:C:247:VAL:H	1.30	0.76
4:C:1484:NAG:HO4	11:C:1486:BMA:C1	2.00	0.74
8:A:1493:MAN:C3	7:A:1494:MAN:C1	2.66	0.73
1:C:207:ASN:HD21	4:C:1483:NAG:C2	1.97	0.72
1:D:471:GLU:HG3	1:D:473:ILE:HG22	1.72	0.71
1:A:93:LEU:H	1:A:240:HIS:CD2	2.09	0.71
1:C:423:LYS:HZ3	10:C:1478:PEG:H42	1.54	0.71
1:D:207:ASN:HD21	5:D:1481:NAG:C2	2.00	0.69
1:C:423:LYS:HZ3	10:C:1478:PEG:H32	1.57	0.68
1:B:233:GLN:HE21	1:B:247:VAL:H	1.42	0.67
9:B:1487:NAG:C1	5:D:1481:NAG:H4	2.24	0.67
1:D:471:GLU:CG	1:D:473:ILE:HG22	2.24	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:1488:NAG:H61	1:C:461:LEU:O	1.95	0.67
1:D:233:GLN:HE21	1:D:247:VAL:H	1.42	0.66
1:B:207:ASN:HD21	5:B:1482:NAG:C2	2.09	0.66
1:A:233:GLN:HE21	1:A:247:VAL:H	1.42	0.66
6:A:1483:MAN:C3	7:A:1485:MAN:C1	2.74	0.65
1:D:93:LEU:H	1:D:240:HIS:CD2	2.07	0.65
1:D:342:ASN:HD22	1:D:344:ASP:H	1.44	0.65
1:C:152:ASN:ND2	5:C:1481:NAG:C2	2.60	0.64
1:C:93:LEU:H	1:C:240:HIS:CD2	2.04	0.64
1:B:405:ASN:HD22	1:B:405:ASN:C	2.02	0.63
1:C:150:HIS:CE1	1:D:471:GLU:H	2.06	0.62
11:C:1487:MAN:H2	1:D:400:ASN:HB3	1.81	0.62
1:B:247:VAL:HG22	1:B:261:ILE:HD12	1.81	0.62
1:D:342:ASN:ND2	1:D:344:ASP:H	1.98	0.61
1:B:207:ASN:HD21	5:B:1482:NAG:H2	1.65	0.61
1:C:465:SER:OG	1:C:467:HIS:HD2	1.84	0.60
1:C:423:LYS:HZ2	10:C:1478:PEG:H32	1.64	0.60
1:D:152:ASN:ND2	5:D:1480:NAG:C2	2.64	0.60
1:B:207:ASN:ND2	5:B:1482:NAG:H2	2.17	0.59
1:C:458:LYS:HE3	12:C:2011:HOH:O	2.03	0.59
1:C:152:ASN:ND2	5:C:1481:NAG:O5	2.26	0.59
8:A:1493:MAN:HO3	7:A:1494:MAN:C1	2.12	0.59
1:C:427:ASN:OD1	10:C:1478:PEG:H22	2.03	0.59
1:A:473:ILE:HG12	12:A:2421:HOH:O	2.03	0.58
1:D:88:ARG:HH11	1:D:88:ARG:CG	2.15	0.58
1:B:400:ASN:O	9:B:1487:NAG:O7	2.21	0.58
1:C:207:ASN:ND2	4:C:1483:NAG:C2	2.62	0.58
8:A:1488:NAG:O5	1:C:462:GLY:HA2	2.05	0.57
1:A:324:LEU:HD21	1:A:362:LEU:HD21	1.85	0.57
6:A:1483:MAN:HO6	7:A:1484:MAN:C1	2.12	0.57
1:B:207:ASN:ND2	5:B:1482:NAG:C1	2.65	0.56
1:C:152:ASN:CG	5:C:1481:NAG:C1	2.68	0.56
3:A:1478:GOL:H32	1:C:113:ARG:NH2	2.22	0.55
1:C:423:LYS:HZ3	10:C:1478:PEG:C4	2.17	0.54
1:B:471:GLU:H	1:D:150:HIS:CE1	2.07	0.54
1:A:342:ASN:HD22	1:A:344:ASP:H	1.54	0.54
1:A:150:HIS:CE1	1:C:471:GLU:H	2.09	0.54
1:A:465:SER:OG	1:A:467:HIS:HD2	1.91	0.54
1:B:240:HIS:HE1	1:B:315:GLU:OE2	1.90	0.54
1:A:353:ASP:OD2	12:A:2301:HOH:O	0.53	0.53
1:D:373:ILE:HD13	7:D:1483:MAN:H61	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:TYR:CG	1:B:235:SER:HA	2.43	0.53
10:C:1478:PEG:H12	12:C:2382:HOH:O	2.08	0.53
1:C:233:GLN:NE2	1:C:247:VAL:H	2.04	0.53
1:A:302:TRP:O	1:A:303:LYS:HE2	2.09	0.53
1:B:342:ASN:HD22	1:B:344:ASP:H	1.56	0.53
1:A:124:ARG:HA	1:A:449:ASN:ND2	2.24	0.53
1:D:92:ASN:HD21	5:D:1479:NAG:C1	2.22	0.52
1:C:240:HIS:HE1	1:C:315:GLU:OE1	1.92	0.52
1:C:102:SER:HB2	1:C:459:LYS:O	2.09	0.52
1:B:152:ASN:ND2	5:B:1480:NAG:C2	2.68	0.52
1:A:467:HIS:H	1:A:467:HIS:CD2	2.28	0.52
1:C:295:LYS:HE3	12:C:2249:HOH:O	2.09	0.52
3:A:1478:GOL:H32	1:C:113:ARG:HH22	1.75	0.51
1:A:93:LEU:N	1:A:240:HIS:HD2	2.02	0.51
11:C:1487:MAN:H2	1:D:400:ASN:CB	2.39	0.51
1:B:207:ASN:ND2	5:B:1482:NAG:C2	2.74	0.51
1:D:94:THR:HG21	5:D:1479:NAG:H81	1.93	0.51
1:D:281:HIS:HD2	1:D:301:ASN:H	1.59	0.51
1:D:240:HIS:HE1	1:D:315:GLU:OE1	1.94	0.50
1:D:88:ARG:CB	1:D:88:ARG:NH1	2.54	0.50
1:B:342:ASN:ND2	1:B:344:ASP:H	2.09	0.50
1:D:340:ASN:ND2	1:D:394:GLN:HE21	2.08	0.50
6:A:1495:NAG:C1	5:B:1482:NAG:C4	2.89	0.50
1:C:127:TYR:CG	1:C:235:SER:HA	2.47	0.49
1:A:423:LYS:HE3	12:A:2381:HOH:O	2.12	0.49
4:C:1484:NAG:C4	11:C:1486:BMA:C1	2.91	0.49
5:C:1482:NAG:H62	1:D:473:ILE:HD13	1.95	0.49
1:A:240:HIS:HE1	1:A:315:GLU:OE2	1.96	0.49
1:B:152:ASN:CG	5:B:1480:NAG:C1	2.72	0.48
1:A:88:ARG:HH11	1:A:194:MET:HA	1.77	0.48
1:A:218:PRO:HG2	10:C:1478:PEG:H11	1.94	0.48
1:D:92:ASN:ND2	5:D:1479:NAG:C1	2.77	0.48
1:A:131:ASP:HB2	1:A:132:PRO:CD	2.44	0.47
1:D:127:TYR:CG	1:D:235:SER:HA	2.49	0.47
1:B:124:ARG:HA	1:B:449:ASN:ND2	2.30	0.47
1:A:407:GLN:NE2	12:A:2363:HOH:O	2.43	0.47
1:D:340:ASN:HD21	1:D:394:GLN:HE21	1.62	0.47
1:A:127:TYR:CG	1:A:235:SER:HA	2.50	0.47
1:C:340:ASN:ND2	12:C:2292:HOH:O	2.47	0.46
1:C:281:HIS:HD2	1:C:301:ASN:H	1.63	0.46
1:C:423:LYS:HZ1	10:C:1478:PEG:H31	1.78	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:247:VAL:HG22	1:C:261:ILE:HD12	1.97	0.46
8:A:1488:NAG:H81	12:A:2148:HOH:O	2.15	0.46
1:D:322:LYS:HB2	1:D:343:CYS:O	2.16	0.45
1:B:324:LEU:HD21	1:B:362:LEU:HD21	1.98	0.45
1:D:373:ILE:HG21	7:D:1483:MAN:H61	2.00	0.44
1:C:467:HIS:HE1	12:C:2421:HOH:O	2.00	0.44
1:A:99:GLU:HB3	1:A:458:LYS:HD3	2.00	0.43
1:A:342:ASN:ND2	1:A:344:ASP:H	2.15	0.43
1:A:113:ARG:NH2	3:B:1478:GOL:H32	2.33	0.43
1:B:465:SER:OG	1:B:467:HIS:HD2	2.01	0.43
1:C:417:ILE:HB	1:C:419:TYR:CZ	2.54	0.42
1:C:228:ASN:HB3	1:C:251:GLY:HA2	2.01	0.42
1:A:195:SER:HB2	1:A:214:TYR:CZ	2.54	0.42
9:B:1487:NAG:O7	5:D:1481:NAG:H61	2.19	0.42
1:C:427:ASN:OD1	10:C:1478:PEG:C2	2.68	0.42
1:D:131:ASP:HB2	1:D:132:PRO:CD	2.49	0.42
1:D:341:GLY:HA2	1:D:347:ILE:HD11	2.01	0.42
1:C:150:HIS:HE1	1:D:471:GLU:N	1.97	0.42
1:C:285:CYS:HB3	1:C:296:CYS:HB3	2.01	0.42
1:D:369:LEU:HG	1:D:386:VAL:HG13	2.02	0.41
1:D:424:GLU:HG2	1:D:424:GLU:O	2.20	0.41
1:C:305:ALA:HB2	1:C:349:GLY:HA2	2.01	0.41
1:D:195:SER:HB2	1:D:214:TYR:CZ	2.55	0.41
1:D:152:ASN:CG	5:D:1480:NAG:C1	2.77	0.41
1:D:342:ASN:HD22	1:D:342:ASN:C	2.24	0.41
1:B:471:GLU:HB2	1:B:474:TYR:CD1	2.56	0.41
1:D:417:ILE:HB	1:D:419:TYR:CZ	2.56	0.41
1:D:471:GLU:OE2	1:D:473:ILE:HG22	2.21	0.41
1:A:150:HIS:HE1	1:C:471:GLU:N	1.99	0.41
3:C:1479:GOL:H32	1:D:113:ARG:NH2	2.35	0.41
1:D:124:ARG:HA	1:D:449:ASN:ND2	2.35	0.41
8:A:1493:MAN:HO6	7:C:1485:MAN:C1	2.23	0.40
1:B:342:ASN:C	1:B:342:ASN:HD22	2.24	0.40
1:B:417:ILE:HG21	1:B:426:PHE:HB3	2.03	0.40
1:D:207:ASN:ND2	5:D:1481:NAG:C2	2.70	0.40
1:B:467:HIS:CD2	1:B:467:HIS:H	2.38	0.40
1:C:325:CYS:O	1:C:393:ILE:HA	2.21	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	387/389 (100%)	375 (97%)	12 (3%)	0	100	100
1	B	387/389 (100%)	375 (97%)	12 (3%)	0	100	100
1	C	387/389 (100%)	374 (97%)	13 (3%)	0	100	100
1	D	387/389 (100%)	373 (96%)	14 (4%)	0	100	100
All	All	1548/1556 (100%)	1497 (97%)	51 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	331/331 (100%)	327 (99%)	4 (1%)	78	69
1	B	331/331 (100%)	325 (98%)	6 (2%)	66	52
1	C	331/331 (100%)	327 (99%)	4 (1%)	78	69
1	D	331/331 (100%)	323 (98%)	8 (2%)	57	39
All	All	1324/1324 (100%)	1302 (98%)	22 (2%)	68	54

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	LEU
1	A	342	ASN

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Mol	Chain	Res	Type
1	A	445	LEU
1	A	464	TRP
1	B	88	ARG
1	B	128	LEU
1	B	142	GLN
1	B	342	ASN
1	B	405	ASN
1	B	464	TRP
1	C	128	LEU
1	C	324	LEU
1	C	445	LEU
1	C	464	TRP
1	D	88	ARG
1	D	128	LEU
1	D	342	ASN
1	D	386	VAL
1	D	424	GLU
1	D	445	LEU
1	D	460	ARG
1	D	464	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (52) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	HIS
1	A	233	GLN
1	A	240	HIS
1	A	340	ASN
1	A	342	ASN
1	A	400	ASN
1	A	406	ASN
1	A	407	GLN
1	A	408	ASN
1	A	422	ASN
1	A	427	ASN
1	A	449	ASN
1	A	467	HIS
1	B	142	GLN
1	B	152	ASN
1	B	207	ASN
1	B	233	GLN
1	B	240	HIS

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Mol	Chain	Res	Type
1	B	340	ASN
1	B	342	ASN
1	B	405	ASN
1	B	406	ASN
1	B	408	ASN
1	B	422	ASN
1	B	427	ASN
1	B	449	ASN
1	B	467	HIS
1	C	150	HIS
1	C	152	ASN
1	C	207	ASN
1	C	233	GLN
1	C	240	HIS
1	C	281	HIS
1	C	340	ASN
1	C	394	GLN
1	C	406	ASN
1	C	408	ASN
1	C	422	ASN
1	C	449	ASN
1	C	467	HIS
1	D	150	HIS
1	D	152	ASN
1	D	207	ASN
1	D	233	GLN
1	D	240	HIS
1	D	281	HIS
1	D	340	ASN
1	D	342	ASN
1	D	406	ASN
1	D	422	ASN
1	D	427	ASN
1	D	449	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 ⓘ

19 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$
4	NAG	A	1479	1,4	14,14,15	0.62	0	15,19,21	0.86	0
4	NAG	A	1480	4	14,14,15	0.53	0	15,19,21	0.74	1 (6%)
6	MAN	A	1483	7,6	11,11,12	0.56	0	14,15,17	1.50	2 (14%)
6	BMA	A	1486	6	11,11,12	0.26	0	14,15,17	1.29	2 (14%)
6	MAN	A	1487	6	11,11,12	0.40	0	14,15,17	0.89	0
8	NAG	A	1488	1,8	14,14,15	0.75	0	15,19,21	2.47	5 (33%)
8	NAG	A	1489	8	14,14,15	0.60	0	15,19,21	1.51	2 (13%)
8	BMA	A	1490	8	11,11,12	0.54	0	14,15,17	1.05	1 (7%)
8	MAN	A	1491	8	11,11,12	0.50	0	14,15,17	0.87	1 (7%)
8	MAN	A	1492	8	11,11,12	0.42	0	14,15,17	0.96	1 (7%)
8	MAN	A	1493	8,7	11,11,12	0.40	0	14,15,17	1.38	2 (14%)
6	NAG	A	1495	5,6	14,14,15	0.57	0	15,19,21	0.62	0
9	MAN	B	1483	9,7	11,11,12	0.43	0	14,15,17	1.08	1 (7%)
9	BMA	B	1486	9	11,11,12	0.40	0	14,15,17	1.24	1 (7%)
9	NAG	B	1487	9	14,14,15	0.65	0	15,19,21	2.50	5 (33%)
4	NAG	C	1483	1,4	14,14,15	0.59	0	15,19,21	1.67	3 (20%)
4	NAG	C	1484	4	14,14,15	0.58	0	15,19,21	0.54	0
11	BMA	C	1486	11	11,11,12	0.42	0	14,15,17	1.11	1 (7%)
11	MAN	C	1487	11	11,11,12	0.47	0	14,15,17	1.69	2 (14%)

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	A	1479	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	1480	4	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	MAN	A	1483	7,6	-	0/2/19/22	0/1/1/1
6	BMA	A	1486	6	-	0/2/19/22	0/1/1/1
6	MAN	A	1487	6	-	0/2/19/22	0/1/1/1
8	NAG	A	1488	1,8	-	0/6/23/26	0/1/1/1
8	NAG	A	1489	8	-	0/6/23/26	0/1/1/1
8	BMA	A	1490	8	-	0/2/19/22	0/1/1/1
8	MAN	A	1491	8	-	0/2/19/22	0/1/1/1
8	MAN	A	1492	8	-	0/2/19/22	0/1/1/1
8	MAN	A	1493	8,7	-	0/2/19/22	0/1/1/1
6	NAG	A	1495	5,6	-	0/6/23/26	0/1/1/1
9	MAN	B	1483	9,7	-	0/2/19/22	0/1/1/1
9	BMA	B	1486	9	-	0/2/19/22	0/1/1/1
9	NAG	B	1487	9	-	0/6/23/26	0/1/1/1
4	NAG	C	1483	1,4	-	0/6/23/26	0/1/1/1
4	NAG	C	1484	4	-	0/6/23/26	0/1/1/1
11	BMA	C	1486	11	-	0/2/19/22	0/1/1/1
11	MAN	C	1487	11	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1488	NAG	C3-C4-C5	-6.14	99.49	110.20
9	B	1487	NAG	C4-C3-C2	-5.33	102.94	111.23
4	C	1483	NAG	C1-O5-C5	-4.53	106.50	112.25
8	A	1488	NAG	C1-O5-C5	-3.98	107.20	112.25
11	C	1487	MAN	C2-C3-C4	-3.51	105.07	111.04
6	A	1483	MAN	O6-C6-C5	-3.07	101.18	111.33
8	A	1493	MAN	O6-C6-C5	-2.78	102.14	111.33
4	C	1483	NAG	C2-N2-C7	-2.71	119.55	123.04
11	C	1486	BMA	C1-C2-C3	-2.58	106.49	109.54
8	A	1488	NAG	O6-C6-C5	-2.48	103.13	111.33
9	B	1483	MAN	O6-C6-C5	-2.42	103.33	111.33
8	A	1489	NAG	O4-C4-C5	-2.23	103.33	109.24
6	A	1483	MAN	O5-C1-C2	-2.21	107.26	110.86
9	B	1487	NAG	O4-C4-C5	-2.06	103.79	109.24
8	A	1491	MAN	C1-O5-C5	2.06	114.86	112.25
4	A	1480	NAG	C1-O5-C5	2.14	114.97	112.25
6	A	1486	BMA	C1-C2-C3	2.22	112.16	109.54
9	B	1487	NAG	C1-O5-C5	2.29	115.16	112.25
8	A	1492	MAN	C1-O5-C5	2.34	115.21	112.25
4	C	1483	NAG	C3-C4-C5	2.35	114.29	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1488	NAG	O4-C4-C3	2.75	116.53	110.34
8	A	1493	MAN	C1-O5-C5	2.91	115.94	112.25
11	C	1487	MAN	C1-O5-C5	3.12	116.21	112.25
8	A	1490	BMA	C1-O5-C5	3.32	116.46	112.25
9	B	1486	BMA	C1-O5-C5	3.64	116.87	112.25
8	A	1488	NAG	O5-C5-C6	3.68	115.32	107.35
6	A	1486	BMA	C1-O5-C5	3.80	117.08	112.25
8	A	1489	NAG	C1-O5-C5	4.44	117.88	112.25
9	B	1487	NAG	C3-C2-N2	4.51	121.36	110.56
9	B	1487	NAG	C2-N2-C7	4.87	129.29	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1483	MAN	2	0
8	A	1488	NAG	3	0
8	A	1493	MAN	3	0
6	A	1495	NAG	2	0
9	B	1487	NAG	5	0
4	C	1483	NAG	3	0
4	C	1484	NAG	3	0
11	C	1486	BMA	3	0
11	C	1487	MAN	2	0

5.6 Ligand geometry

Of 28 ligands modelled in this entry, 4 are monoatomic - leaving 24 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
3	GOL	A	1478	-	5,5,5	1.17	0	5,5,5	1.81	1 (20%)
5	NAG	A	1481	1	14,14,15	0.62	0	15,19,21	0.81	0
5	NAG	A	1482	-	14,14,15	0.52	0	15,19,21	1.01	1 (6%)
7	MAN	A	1484	6	11,11,12	0.36	0	14,15,17	1.53	2 (14%)
7	MAN	A	1485	6	11,11,12	0.50	0	14,15,17	1.44	2 (14%)
7	MAN	A	1494	8	11,11,12	0.58	0	14,15,17	1.91	4 (28%)
3	GOL	B	1478	-	5,5,5	1.69	1 (20%)	5,5,5	2.13	3 (60%)
5	NAG	B	1479	-	14,14,15	0.43	0	15,19,21	1.20	2 (13%)
5	NAG	B	1480	1	14,14,15	0.68	0	15,19,21	1.09	1 (6%)
5	NAG	B	1481	-	14,14,15	0.50	0	15,19,21	1.03	1 (6%)
5	NAG	B	1482	6	14,14,15	0.55	0	15,19,21	1.40	3 (20%)
7	MAN	B	1484	9	11,11,12	0.37	0	14,15,17	1.22	2 (14%)
7	MAN	B	1485	-	11,11,12	0.50	0	14,15,17	1.95	3 (21%)
10	PEG	C	1478	-	6,6,6	0.26	0	5,5,5	5.99	4 (80%)
3	GOL	C	1479	-	5,5,5	1.59	1 (20%)	5,5,5	2.11	2 (40%)
5	NAG	C	1480	-	14,14,15	0.53	0	15,19,21	1.43	3 (20%)
5	NAG	C	1481	1	14,14,15	0.64	0	15,19,21	1.21	1 (6%)
5	NAG	C	1482	-	14,14,15	0.48	0	15,19,21	0.74	0
7	MAN	C	1485	8	11,11,12	0.42	0	14,15,17	1.66	2 (14%)
3	GOL	D	1478	-	5,5,5	1.35	0	5,5,5	1.87	1 (20%)
5	NAG	D	1479	-	14,14,15	0.44	0	15,19,21	1.04	0
5	NAG	D	1480	1	14,14,15	0.58	0	15,19,21	1.30	1 (6%)
5	NAG	D	1481	1	14,14,15	0.60	0	15,19,21	1.81	3 (20%)
7	MAN	D	1483	-	11,11,12	0.58	0	14,15,17	1.30	2 (14%)

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	GOL	A	1478	-	-	0/4/4/4	0/0/0/0
5	NAG	A	1481	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1482	-	-	0/6/23/26	0/1/1/1
7	MAN	A	1484	6	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	MAN	A	1485	6	-	0/2/19/22	0/1/1/1
7	MAN	A	1494	8	-	0/2/19/22	0/1/1/1
3	GOL	B	1478	-	-	0/4/4/4	0/0/0/0
5	NAG	B	1479	-	-	0/6/23/26	0/1/1/1
5	NAG	B	1480	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1481	-	-	0/6/23/26	0/1/1/1
5	NAG	B	1482	6	-	0/6/23/26	0/1/1/1
7	MAN	B	1484	9	-	0/2/19/22	0/1/1/1
7	MAN	B	1485	-	-	0/2/19/22	0/1/1/1
10	PEG	C	1478	-	-	0/4/4/4	0/0/0/0
3	GOL	C	1479	-	-	0/4/4/4	0/0/0/0
5	NAG	C	1480	-	-	0/6/23/26	0/1/1/1
5	NAG	C	1481	1	-	0/6/23/26	0/1/1/1
5	NAG	C	1482	-	-	0/6/23/26	0/1/1/1
7	MAN	C	1485	8	-	0/2/19/22	0/1/1/1
3	GOL	D	1478	-	-	0/4/4/4	0/0/0/0
5	NAG	D	1479	-	-	0/6/23/26	0/1/1/1
5	NAG	D	1480	1	-	0/6/23/26	0/1/1/1
5	NAG	D	1481	1	-	0/6/23/26	0/1/1/1
7	MAN	D	1483	-	-	0/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1479	GOL	O3-C3	2.29	1.52	1.42
3	B	1478	GOL	O3-C3	2.65	1.53	1.42

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	C	1478	PEG	C3-O2-C2	-6.20	86.66	113.31
5	D	1481	NAG	C1-O5-C5	-4.67	106.32	112.25
5	D	1480	NAG	C1-O5-C5	-3.80	107.42	112.25
5	C	1481	NAG	C1-O5-C5	-2.73	108.78	112.25
7	A	1494	MAN	O2-C2-C1	-2.61	103.98	109.21
5	B	1482	NAG	C4-C3-C2	-2.37	107.55	111.23
5	B	1481	NAG	C2-N2-C7	-2.33	120.04	123.04
5	B	1480	NAG	C1-O5-C5	-2.25	109.39	112.25
5	D	1481	NAG	C2-N2-C7	-2.21	120.19	123.04
5	B	1482	NAG	C2-N2-C7	-2.17	120.25	123.04
7	A	1484	MAN	O2-C2-C1	-2.13	104.93	109.21
5	C	1480	NAG	C4-C3-C2	-2.12	107.94	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	1484	MAN	O2-C2-C1	-2.05	105.10	109.21
3	B	1478	GOL	O1-C1-C2	2.03	120.01	110.18
3	B	1478	GOL	C3-C2-C1	2.12	119.45	111.12
7	D	1483	MAN	C3-C4-C5	2.39	114.36	110.20
5	B	1482	NAG	O5-C5-C6	2.40	112.55	107.35
3	C	1479	GOL	O1-C1-C2	2.46	122.13	110.18
7	A	1494	MAN	O5-C1-C2	2.49	114.89	110.86
5	B	1479	NAG	C3-C4-C5	2.51	114.57	110.20
7	C	1485	MAN	O5-C1-C2	2.53	114.96	110.86
7	A	1485	MAN	O5-C1-C2	2.53	114.96	110.86
5	C	1480	NAG	C3-C4-C5	2.65	114.82	110.20
5	A	1482	NAG	C3-C4-C5	2.81	115.10	110.20
5	C	1480	NAG	C1-O5-C5	2.93	115.97	112.25
5	B	1479	NAG	C1-O5-C5	3.07	116.15	112.25
10	C	1478	PEG	O2-C3-C4	3.17	125.01	110.43
3	A	1478	GOL	O3-C3-C2	3.27	126.05	110.18
7	B	1484	MAN	C1-O5-C5	3.35	116.50	112.25
3	C	1479	GOL	O3-C3-C2	3.38	126.59	110.18
7	B	1485	MAN	O5-C1-C2	3.39	116.35	110.86
7	B	1485	MAN	C1-C2-C3	3.45	113.62	109.54
7	A	1494	MAN	C1-C2-C3	3.45	113.62	109.54
7	D	1483	MAN	C1-C2-C3	3.56	113.75	109.54
3	D	1478	GOL	O3-C3-C2	3.59	127.62	110.18
3	B	1478	GOL	O3-C3-C2	3.64	127.82	110.18
5	D	1481	NAG	C3-C4-C5	3.89	116.99	110.20
7	A	1485	MAN	C1-O5-C5	4.09	117.44	112.25
7	A	1484	MAN	C1-O5-C5	4.22	117.61	112.25
7	A	1494	MAN	C1-O5-C5	4.68	118.19	112.25
7	C	1485	MAN	C1-O5-C5	5.00	118.59	112.25
7	B	1485	MAN	C1-O5-C5	5.21	118.86	112.25
10	C	1478	PEG	O2-C2-C1	7.40	144.49	110.43
10	C	1478	PEG	O1-C1-C2	8.52	164.65	112.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

16 monomers are involved in 56 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1478	GOL	2	0
7	A	1484	MAN	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1485	MAN	1	0
7	A	1494	MAN	2	0
3	B	1478	GOL	1	0
5	B	1480	NAG	3	0
5	B	1482	NAG	8	0
10	C	1478	PEG	16	0
3	C	1479	GOL	1	0
5	C	1481	NAG	4	0
5	C	1482	NAG	1	0
7	C	1485	MAN	1	0
5	D	1479	NAG	3	0
5	D	1480	NAG	3	0
5	D	1481	NAG	7	0
7	D	1483	MAN	2	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 [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	389/389 (100%)	-0.45	4 (1%) 84 83	7, 12, 20, 31	0
1	B	389/389 (100%)	-0.47	1 (0%) 94 93	7, 12, 18, 37	0
1	C	389/389 (100%)	-0.49	6 (1%) 76 74	7, 11, 19, 33	0
1	D	389/389 (100%)	-0.45	3 (0%) 87 86	7, 12, 18, 33	0
All	All	1556/1556 (100%)	-0.46	14 (0%) 85 85	7, 12, 19, 37	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	88	ARG	4.3
1	C	350	GLY	3.6
1	C	349	GLY	3.3
1	A	88	ARG	3.2
1	C	88	ARG	2.9
1	D	88	ARG	2.9
1	A	460	ARG	2.8
1	C	256	ARG	2.7
1	D	424	GLU	2.6
1	D	460	ARG	2.3
1	A	256	ARG	2.2
1	C	351	SER	2.1
1	C	424	GLU	2.1
1	A	133	GLN	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 ⓘ

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
4	NAG	C	1483	14/15	0.86	0.24	12.69	25,29,32,34	0
9	NAG	B	1487	14/15	0.69	0.26	11.35	26,35,39,43	0
6	MAN	A	1483	11/12	0.90	0.20	7.38	22,24,24,27	0
8	MAN	A	1493	11/12	0.95	0.14	5.41	14,15,16,19	0
8	NAG	A	1488	14/15	0.81	0.32	4.24	26,31,35,38	0
6	NAG	A	1495	14/15	0.80	0.28	2.32	20,24,27,30	0
4	NAG	C	1484	14/15	0.81	0.21	1.59	18,24,29,33	0
8	NAG	A	1489	14/15	0.92	0.15	1.00	14,20,24,25	0
9	BMA	B	1486	11/12	0.85	0.24	-	35,36,37,38	0
6	BMA	A	1486	11/12	0.94	0.21	-	21,25,26,31	0
4	NAG	A	1480	14/15	0.68	0.45	-	39,43,45,45	0
8	MAN	A	1492	11/12	0.84	0.41	-	38,42,44,44	0
11	MAN	C	1487	11/12	0.60	0.42	-	45,49,51,51	0
8	MAN	A	1491	11/12	0.94	0.13	-	24,25,27,27	0
4	NAG	A	1479	14/15	0.90	0.16	-	20,22,26,32	0
8	BMA	A	1490	11/12	0.94	0.13	-	16,19,19,21	0
9	MAN	B	1483	11/12	0.82	0.26	-	35,37,37,37	0
11	BMA	C	1486	11/12	0.85	0.28	-	32,35,36,40	0
6	MAN	A	1487	11/12	0.76	0.44	-	45,48,49,50	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(Å ²)	Q<0.9
7	MAN	B	1485	11/12	0.74	0.33	12.17	44,47,48,49	0
3	GOL	B	1478	6/6	0.67	0.28	9.34	12,25,26,28	0
3	GOL	D	1478	6/6	0.73	0.29	7.34	17,29,30,30	0
3	GOL	C	1479	6/6	0.76	0.26	7.12	13,26,28,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
7	MAN	A	1485	11/12	0.85	0.18	6.45	24,27,28,29	0
5	NAG	D	1481	14/15	0.82	0.21	6.34	23,26,27,27	0
3	GOL	A	1478	6/6	0.69	0.24	6.32	19,29,30,31	0
10	PEG	C	1478	7/7	0.67	0.25	6.31	20,24,28,29	0
2	CA	C	1477	1/1	0.98	0.17	5.29	19,19,19,19	0
2	CA	D	1477	1/1	0.99	0.13	3.76	19,19,19,19	0
7	MAN	A	1494	11/12	0.92	0.14	3.57	19,20,21,21	0
7	MAN	C	1485	11/12	0.85	0.19	2.99	30,33,36,38	0
5	NAG	B	1482	14/15	0.73	0.30	2.01	42,44,45,45	0
2	CA	B	1477	1/1	0.99	0.09	1.68	19,19,19,19	0
2	CA	A	1477	1/1	0.99	0.14	1.36	20,20,20,20	0
5	NAG	B	1481	14/15	0.15	0.47	-	95,96,97,97	0
5	NAG	C	1482	14/15	0.42	0.37	-	74,77,77,77	0
5	NAG	A	1482	14/15	0.47	0.49	-	64,67,68,68	0
5	NAG	B	1480	14/15	0.90	0.14	-	16,22,26,29	0
5	NAG	A	1481	14/15	0.89	0.16	-	19,22,27,28	0
5	NAG	C	1481	14/15	0.93	0.13	-	15,18,25,26	0
7	MAN	D	1483	11/12	0.53	0.39	-	67,68,68,68	0
7	MAN	A	1484	11/12	0.81	0.20	-	26,30,35,38	0
5	NAG	C	1480	14/15	0.47	0.37	-	62,63,64,64	0
5	NAG	D	1480	14/15	0.92	0.15	-	17,22,26,26	0
5	NAG	B	1479	14/15	0.31	0.43	-	98,99,99,99	0
7	MAN	B	1484	11/12	0.79	0.29	-	32,33,36,38	0
5	NAG	D	1479	14/15	0.50	0.43	-	105,105,106,106	0

6.5 Other polymers [i](#)

There are no such residues in this entry.