



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 10:10 am GMT

PDB ID : 3HTT
Title : The hemagglutinin structure of an avian H1N1 influenza A virus in complex with 2,3-sialyllactose
Authors : Wang, G.; Li, A.; Zhang, Q.; Wu, C.; Zhang, R.; Cai, Q.; Song, W.; Yuen, K.-Y.
Deposited on : 2009-06-12
Resolution : 2.95 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

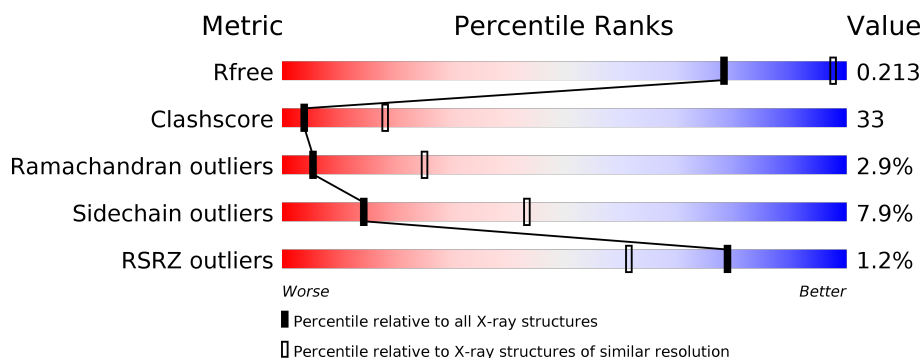
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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}	100719	2395 (3.00-2.92)
Clashscore	112137	2773 (3.00-2.92)
Ramachandran outliers	110173	2680 (3.00-2.92)
Sidechain outliers	110143	2683 (3.00-2.92)
RSRZ outliers	101464	2421 (3.00-2.92)

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	324	<div> <div>2%</div> <div> <div></div> <div>46%</div> <div>47%</div> <div>6%</div> </div> </div>
2	B	160	<div> <div>%</div> <div> <div></div> <div>58%</div> <div>37%</div> <div>6%</div> </div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4203 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 Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2530	1596	432	491	11			

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	160	Total	C	N	O	S	0	0	0
			1286	803	220	256	7			

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

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

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	3	Total	C	N	O	0	0
			43	23	1	19		

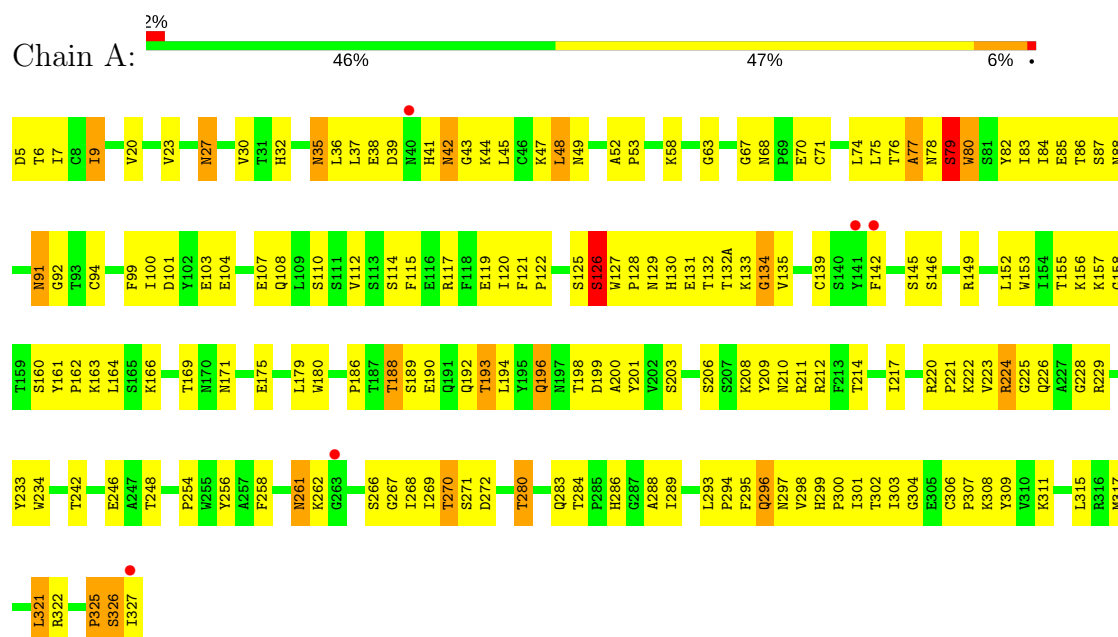
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	184	Total	O	0	0
			184	184		
6	B	104	Total	O	0	0
			104	104		

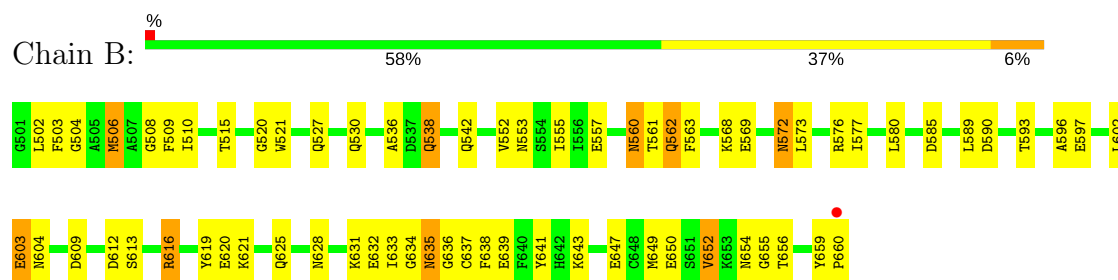
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Hemagglutinin



• Molecule 2: Hemagglutinin HA2 chain



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	198.53Å 198.53Å 198.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.79 – 2.95 46.79 – 2.95	Depositor EDS
% Data completeness (in resolution range)	89.7 (46.79-2.95) 100.0 (46.79-2.95)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.88 (at 2.96Å)	Xtriage
Refinement program	CNS 1.21	Depositor
R, R_{free}	0.202 , 0.247 0.209 , 0.213	Depositor DCC
R_{free} test set	1375 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	54.7	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 57.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.024 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4203	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: GAL, SIA, GLC, NAG, NDG

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	0.35	0/2594	0.62	0/3534
2	B	0.40	0/1312	0.62	0/1766
All	All	0.37	0/3906	0.62	0/5300

There are no bond length outliers.

There are no bond angle outliers.

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	2530	0	2444	193	1
2	B	1286	0	1197	81	0
3	A	28	0	25	7	0
4	A	28	0	26	2	0
5	A	43	0	37	4	0
6	A	184	0	0	43	0
6	B	104	0	0	12	1
All	All	4203	0	3729	255	1

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 33.

All (255) 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:78:ASN:HA	6:A:492:HOH:O	1.57	1.02
1:A:156:LYS:HD3	1:A:196:GLN:HB2	1.50	0.93
6:A:437:HOH:O	2:B:589:LEU:HB3	1.67	0.93
1:A:122:PRO:O	1:A:126:SER:HB2	1.70	0.91
1:A:110:SER:HB2	6:A:411:HOH:O	1.73	0.89
1:A:134:GLY:HA3	1:A:153:TRP:HB3	1.56	0.88
2:B:538:GLN:HG3	6:B:214:HOH:O	1.80	0.80
1:A:42:ASN:HD21	1:A:288:ALA:H	1.31	0.79
1:A:175:GLU:HG2	6:A:343:HOH:O	1.82	0.78
2:B:503:PHE:O	2:B:616:ARG:HD3	1.84	0.78
1:A:86:THR:HG22	1:A:88:ASN:H	1.48	0.77
1:A:9:ILE:HG21	2:B:619:TYR:HA	1.67	0.76
1:A:190:GLU:HA	1:A:193:THR:CG2	2.17	0.75
1:A:157:LYS:HA	6:A:362:HOH:O	1.85	0.75
1:A:246:GLU:HG2	6:A:424:HOH:O	1.88	0.73
1:A:284:THR:HG22	1:A:302:THR:HG22	1.69	0.73
1:A:152:LEU:HG	6:A:350:HOH:O	1.88	0.72
1:A:115:PHE:HB3	6:A:492:HOH:O	1.90	0.72
1:A:166:LYS:HE2	6:A:464:HOH:O	1.89	0.71
1:A:134:GLY:CA	1:A:153:TRP:HB3	2.20	0.71
1:A:299:HIS:HD2	1:A:301:ILE:H	1.38	0.71
2:B:542:GLN:HG2	6:B:214:HOH:O	1.90	0.70
1:A:119:GLU:OE2	1:A:122:PRO:HA	1.92	0.70
1:A:156:LYS:CD	1:A:196:GLN:HB2	2.21	0.69
1:A:169:THR:HG22	1:A:242:THR:OG1	1.92	0.69
2:B:633:ILE:O	2:B:635:ASN:N	2.21	0.69
1:A:70:GLU:HG3	3:A:626:NAG:HN2	1.56	0.69
1:A:160:SER:HA	6:A:490:HOH:O	1.93	0.69
1:A:156:LYS:HA	1:A:161:TYR:HD1	1.58	0.69
5:A:3033:SIA:O1B	5:A:3033:SIA:H6	1.92	0.68
2:B:560:ASN:HD22	2:B:561:THR:H	1.42	0.68
1:A:9:ILE:C	1:A:9:ILE:HD13	2.14	0.67
1:A:300:PRO:HG3	1:A:309:TYR:CZ	2.29	0.67
1:A:91:ASN:ND2	6:A:380:HOH:O	2.27	0.67
1:A:189:SER:O	1:A:192:GLN:HG3	1.95	0.67
1:A:149:ARG:HG2	1:A:149:ARG:HH11	1.59	0.67
2:B:625:GLN:HE22	2:B:655:GLY:HA2	1.60	0.66
2:B:635:ASN:HD21	2:B:637:CYS:HB2	1.59	0.66
1:A:226:GLN:HB3	6:A:337:HOH:O	1.95	0.66
2:B:553:ASN:O	2:B:557:GLU:HG2	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LYS:HD3	6:A:446:HOH:O	1.96	0.65
1:A:27:ASN:ND2	4:A:632:NAG:H82	2.11	0.65
1:A:42:ASN:ND2	1:A:288:ALA:H	1.94	0.65
1:A:103:GLU:HB2	6:A:4:HOH:O	1.97	0.65
1:A:194:LEU:HD21	5:A:3033:SIA:C11	2.27	0.64
1:A:280:THR:HG21	6:A:368:HOH:O	1.95	0.64
1:A:198:THR:HG21	6:A:438:HOH:O	1.97	0.64
1:A:114:SER:OG	1:A:261:ASN:HB2	1.98	0.63
1:A:156:LYS:HE3	6:A:489:HOH:O	1.97	0.63
2:B:647:GLU:HG2	6:B:42:HOH:O	1.97	0.63
1:A:112:VAL:HG11	1:A:115:PHE:HB2	1.78	0.63
2:B:502:LEU:HB2	2:B:609:ASP:OD1	1.99	0.63
1:A:289:ILE:CD1	1:A:298:VAL:HG21	2.29	0.62
2:B:560:ASN:HD22	2:B:561:THR:N	1.96	0.62
1:A:309:TYR:HD2	2:B:589:LEU:HD13	1.64	0.62
2:B:530:GLN:NE2	2:B:530:GLN:HA	2.14	0.62
1:A:179:LEU:HD23	1:A:234:TRP:HB3	1.81	0.61
1:A:317:MET:HE3	2:B:552:VAL:HG22	1.82	0.61
1:A:160:SER:HB2	6:A:417:HOH:O	2.01	0.61
1:A:127:TRP:O	1:A:130:HIS:HB2	2.01	0.61
1:A:327:ILE:HD11	6:B:169:HOH:O	2.01	0.61
2:B:628:ASN:HB2	6:B:139:HOH:O	2.00	0.61
1:A:131:GLU:OE1	1:A:132(A):THR:HG22	2.00	0.61
1:A:52:ALA:O	1:A:80:TRP:HA	2.01	0.60
1:A:83:ILE:HB	1:A:268:ILE:HG12	1.83	0.60
1:A:86:THR:HG22	1:A:87:SER:N	2.15	0.60
1:A:6:THR:HG22	2:B:639:GLU:HA	1.82	0.60
1:A:94:CYS:O	1:A:224:ARG:HD3	2.01	0.60
1:A:36:LEU:HA	1:A:293:LEU:HD22	1.82	0.60
1:A:119:GLU:HG3	1:A:256:TYR:CZ	2.36	0.60
1:A:304:GLY:HA2	2:B:563:PHE:CE1	2.37	0.59
1:A:289:ILE:HD11	1:A:298:VAL:HG11	1.83	0.59
1:A:299:HIS:CD2	1:A:301:ILE:H	2.19	0.58
1:A:131:GLU:HB3	1:A:155:THR:O	2.03	0.58
2:B:509:PHE:CD1	2:B:510:ILE:HG13	2.39	0.58
1:A:211:ARG:NH2	6:A:344:HOH:O	2.36	0.58
1:A:32:HIS:HD2	2:B:521:TRP:HE1	1.52	0.58
2:B:660:PRO:HG3	6:B:73:HOH:O	2.02	0.58
1:A:193:THR:HG21	6:A:338:HOH:O	2.04	0.57
1:A:126:SER:HA	6:A:483:HOH:O	2.04	0.57
2:B:603:GLU:OE1	2:B:603:GLU:HA	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:632:GLU:HG2	2:B:638:PHE:CE2	2.40	0.57
2:B:502:LEU:HD13	6:B:187:HOH:O	2.04	0.56
1:A:39:ASP:HB3	1:A:297:ASN:OD1	2.04	0.56
2:B:560:ASN:ND2	2:B:561:THR:N	2.52	0.56
2:B:628:ASN:HD22	2:B:628:ASN:N	2.03	0.56
2:B:633:ILE:O	2:B:635:ASN:ND2	2.38	0.56
1:A:156:LYS:CA	1:A:161:TYR:HD1	2.18	0.56
1:A:45:LEU:HD13	1:A:84:ILE:HD13	1.87	0.56
1:A:149:ARG:HG2	1:A:149:ARG:NH1	2.20	0.56
1:A:189:SER:O	1:A:193:THR:HG22	2.05	0.56
1:A:304:GLY:HA2	2:B:563:PHE:CD1	2.40	0.56
1:A:171:ASN:HB2	6:A:328:HOH:O	2.06	0.55
1:A:317:MET:CE	2:B:552:VAL:HG22	2.36	0.55
1:A:212:ARG:HA	6:A:349:HOH:O	2.06	0.55
1:A:47:LYS:HG3	6:A:355:HOH:O	2.06	0.55
2:B:568:LYS:O	2:B:569:GLU:HB2	2.06	0.55
2:B:654:ASN:O	2:B:656:THR:HG23	2.06	0.55
1:A:311:LYS:HD2	6:A:437:HOH:O	2.06	0.55
1:A:43:GLY:HA2	1:A:286:HIS:O	2.07	0.55
2:B:504:GLY:O	2:B:508:GLY:HA3	2.07	0.55
1:A:131:GLU:CD	1:A:132(A):THR:HG22	2.28	0.54
1:A:294:PRO:HG2	1:A:295:PHE:CD1	2.41	0.54
2:B:593:THR:O	2:B:597:GLU:HB2	2.06	0.54
2:B:530:GLN:HE21	2:B:530:GLN:HA	1.70	0.54
1:A:32:HIS:CD2	2:B:521:TRP:HE1	2.25	0.54
1:A:192:GLN:HB3	1:A:198:THR:CG2	2.38	0.54
1:A:289:ILE:HD11	1:A:298:VAL:HG21	1.88	0.54
1:A:225:GLY:HA3	6:A:428:HOH:O	2.08	0.54
1:A:194:LEU:HD21	5:A:3033:SIA:H111	1.90	0.54
1:A:190:GLU:HA	1:A:193:THR:HG22	1.90	0.53
1:A:162:PRO:HD2	6:A:370:HOH:O	2.06	0.53
1:A:300:PRO:HG3	1:A:309:TYR:CE2	2.43	0.53
1:A:42:ASN:C	1:A:42:ASN:HD22	2.12	0.53
1:A:80:TRP:N	6:A:351:HOH:O	2.42	0.53
1:A:86:THR:HG23	1:A:272:ASP:OD1	2.09	0.53
1:A:86:THR:CG2	1:A:87:SER:N	2.71	0.52
1:A:127:TRP:N	1:A:128:PRO:HD3	2.24	0.52
1:A:135:VAL:HG12	1:A:145:SER:HB3	1.91	0.52
1:A:309:TYR:CD2	2:B:589:LEU:HD13	2.43	0.52
1:A:53:PRO:HB3	1:A:82:TYR:CE2	2.45	0.52
2:B:585:ASP:O	2:B:589:LEU:HG	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:TYR:O	2:B:589:LEU:HD22	2.10	0.52
1:A:75:LEU:HD22	1:A:117:ARG:NH1	2.25	0.51
1:A:86:THR:HG22	1:A:88:ASN:N	2.21	0.51
2:B:577:ILE:HG12	6:B:406:HOH:O	2.09	0.51
1:A:58:LYS:HA	1:A:88:ASN:O	2.10	0.51
1:A:71:CYS:HB3	1:A:74:LEU:HD12	1.92	0.51
1:A:308:LYS:NZ	2:B:561:THR:HG22	2.25	0.51
1:A:35:ASN:HD22	1:A:36:LEU:H	1.59	0.51
1:A:188:THR:N	1:A:217:ILE:HG21	2.26	0.51
2:B:635:ASN:C	2:B:635:ASN:HD22	2.13	0.51
1:A:203:SER:OG	1:A:246:GLU:HG2	2.11	0.50
1:A:289:ILE:HD12	1:A:298:VAL:HG21	1.94	0.50
1:A:269:ILE:HG12	1:A:303:ILE:HD12	1.92	0.50
1:A:325:PRO:O	1:A:326:SER:HB3	2.12	0.50
1:A:169:THR:HG23	1:A:242:THR:HG23	1.93	0.50
1:A:186:PRO:HD2	1:A:190:GLU:OE2	2.11	0.50
1:A:296:GLN:HE22	1:A:309:TYR:HB2	1.77	0.50
2:B:625:GLN:NE2	2:B:655:GLY:HA2	2.26	0.50
1:A:91:ASN:OD1	3:A:626:NAG:N2	2.44	0.50
1:A:208:LYS:O	1:A:208:LYS:CG	2.60	0.49
1:A:224:ARG:NH2	3:A:626:NAG:O3	2.45	0.49
2:B:530:GLN:HE21	2:B:530:GLN:CA	2.22	0.49
1:A:164:LEU:C	1:A:164:LEU:HD12	2.33	0.49
2:B:621:LYS:HD3	6:B:201:HOH:O	2.12	0.49
1:A:104:GLU:HB2	6:A:329:HOH:O	2.12	0.48
1:A:85:GLU:O	1:A:270:THR:HA	2.14	0.48
1:A:79:SER:N	6:A:334:HOH:O	2.42	0.48
2:B:572:ASN:HD22	2:B:573:LEU:N	2.11	0.48
1:A:58:LYS:HE2	6:A:380:HOH:O	2.12	0.48
1:A:67:GLY:O	1:A:149:ARG:HG3	2.13	0.48
1:A:23:VAL:CG2	2:B:602:LEU:HD23	2.44	0.48
1:A:212:ARG:CA	6:A:349:HOH:O	2.62	0.48
1:A:133:LYS:O	1:A:135:VAL:N	2.47	0.47
1:A:315:LEU:CD1	2:B:596:ALA:HB1	2.44	0.47
1:A:128:PRO:O	1:A:129:ASN:CG	2.53	0.47
1:A:135:VAL:HG13	1:A:146:SER:CA	2.44	0.47
1:A:156:LYS:HA	1:A:161:TYR:CD1	2.44	0.47
1:A:299:HIS:CD2	1:A:300:PRO:HD2	2.49	0.47
1:A:63:GLY:CA	1:A:92:GLY:HA2	2.44	0.47
2:B:631:LYS:HE3	6:B:136:HOH:O	2.14	0.47
1:A:37:LEU:HB2	1:A:315:LEU:HB2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:ARG:HH11	1:A:322:ARG:HG3	1.80	0.47
2:B:628:ASN:N	2:B:628:ASN:ND2	2.63	0.47
1:A:129:ASN:ND2	6:A:450:HOH:O	2.48	0.47
1:A:164:LEU:HA	6:A:433:HOH:O	2.14	0.47
2:B:569:GLU:N	6:B:239:HOH:O	2.27	0.47
1:A:23:VAL:HG21	2:B:602:LEU:HD23	1.97	0.47
2:B:628:ASN:O	2:B:641:TYR:HD1	1.98	0.47
1:A:122:PRO:O	1:A:126:SER:CB	2.55	0.47
1:A:220:ARG:HB2	1:A:221:PRO:HD2	1.96	0.46
1:A:101:ASP:HA	6:A:329:HOH:O	2.15	0.46
1:A:206:SER:HB3	1:A:209:TYR:CB	2.46	0.46
1:A:7:ILE:O	1:A:7:ILE:HG23	2.16	0.46
2:B:635:ASN:ND2	2:B:637:CYS:H	2.13	0.46
1:A:321:LEU:HD23	1:A:321:LEU:N	2.30	0.46
2:B:562:GLN:OE1	2:B:562:GLN:N	2.49	0.45
1:A:223:VAL:O	1:A:224:ARG:C	2.54	0.45
1:A:157:LYS:HE3	6:A:494:HOH:O	2.15	0.45
1:A:35:ASN:HD22	1:A:36:LEU:N	2.13	0.45
1:A:164:LEU:O	1:A:164:LEU:HD12	2.17	0.45
1:A:20:VAL:HB	2:B:604:ASN:ND2	2.32	0.45
1:A:41:HIS:HB3	1:A:289:ILE:HD13	1.99	0.45
1:A:206:SER:HB3	1:A:209:TYR:HB3	1.99	0.44
1:A:222:LYS:HB3	1:A:225:GLY:O	2.17	0.44
1:A:100:ILE:HG13	1:A:233:TYR:CE2	2.52	0.44
2:B:572:ASN:C	2:B:572:ASN:HD22	2.20	0.44
2:B:659:TYR:CG	2:B:659:TYR:O	2.70	0.44
1:A:163:LYS:HE3	6:A:468:HOH:O	2.17	0.44
1:A:175:GLU:OE1	1:A:262:LYS:HE2	2.18	0.44
1:A:190:GLU:HA	1:A:193:THR:HG23	1.96	0.44
1:A:266:SER:OG	1:A:267:GLY:N	2.50	0.44
1:A:228:GLY:O	1:A:229:ARG:HD3	2.18	0.44
1:A:315:LEU:HD13	2:B:596:ALA:HB1	1.99	0.44
1:A:317:MET:HE1	2:B:552:VAL:HG13	1.98	0.44
1:A:180:TRP:HB3	1:A:254:PRO:HG3	2.00	0.44
1:A:35:ASN:ND2	6:A:341:HOH:O	2.48	0.44
1:A:35:ASN:ND2	1:A:37:LEU:H	2.15	0.43
2:B:647:GLU:O	2:B:650:GLU:HB3	2.18	0.43
2:B:649:MET:O	2:B:652:VAL:HG13	2.17	0.43
1:A:311:LYS:NZ	2:B:589:LEU:HB2	2.33	0.43
1:A:42:ASN:O	1:A:44:LYS:HG3	2.19	0.43
1:A:201:TYR:CE2	1:A:248:THR:HG23	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:616:ARG:HH11	2:B:616:ARG:HG2	1.83	0.43
1:A:94:CYS:HB3	3:A:626:NAG:O7	2.19	0.43
1:A:9:ILE:CD1	1:A:9:ILE:C	2.87	0.43
2:B:572:ASN:ND2	6:B:267:HOH:O	2.51	0.43
1:A:157:LYS:HD3	1:A:158:GLY:N	2.34	0.43
1:A:208:LYS:O	1:A:208:LYS:HG2	2.19	0.43
2:B:520:GLY:HA3	2:B:536:ALA:HB1	2.00	0.43
1:A:311:LYS:HZ1	2:B:589:LEU:HB2	1.82	0.43
1:A:41:HIS:HB3	1:A:289:ILE:CD1	2.49	0.43
3:A:627:NDG:H5	3:A:627:NDG:HA	1.84	0.43
1:A:317:MET:SD	2:B:555:ILE:HD12	2.59	0.43
1:A:169:THR:CG2	1:A:242:THR:OG1	2.65	0.43
1:A:70:GLU:HG3	3:A:626:NAG:N2	2.29	0.42
1:A:68:ASN:HB3	1:A:71:CYS:SG	2.59	0.42
1:A:76:THR:O	1:A:77:ALA:C	2.58	0.42
1:A:79:SER:O	1:A:112:VAL:O	2.37	0.42
1:A:196:GLN:HG2	6:A:490:HOH:O	2.18	0.42
1:A:164:LEU:O	1:A:246:GLU:HA	2.19	0.42
1:A:194:LEU:CD2	5:A:3033:SIA:H111	2.49	0.42
1:A:63:GLY:HA3	1:A:92:GLY:HA2	2.00	0.42
1:A:30:VAL:HG12	1:A:32:HIS:H	1.85	0.42
2:B:635:ASN:H	2:B:635:ASN:HD22	1.66	0.42
1:A:306:CYS:HA	1:A:307:PRO:HD3	1.84	0.42
2:B:635:ASN:HD22	2:B:637:CYS:H	1.68	0.42
3:A:627:NDG:H8C3	6:A:408:HOH:O	2.19	0.42
1:A:117:ARG:HG3	1:A:258:PHE:CE2	2.55	0.41
1:A:200:ALA:HA	1:A:248:THR:OG1	2.20	0.41
1:A:317:MET:CE	2:B:552:VAL:HG13	2.50	0.41
1:A:135:VAL:HG13	1:A:146:SER:N	2.35	0.41
1:A:48:LEU:O	1:A:49:ASN:HB2	2.18	0.41
2:B:506:MET:HE2	2:B:612:ASP:N	2.35	0.41
1:A:107:GLU:CD	6:A:367:HOH:O	2.59	0.41
1:A:86:THR:C	1:A:88:ASN:H	2.24	0.41
2:B:560:ASN:ND2	2:B:561:THR:H	2.12	0.41
1:A:112:VAL:CG1	1:A:115:PHE:HB2	2.49	0.41
4:A:632:NAG:H5	6:A:353:HOH:O	2.20	0.41
1:A:211:ARG:HG2	6:A:1:HOH:O	2.20	0.41
2:B:503:PHE:CE1	2:B:613:SER:HB2	2.56	0.41
2:B:553:ASN:O	2:B:557:GLU:N	2.46	0.41
2:B:643:LYS:HA	2:B:643:LYS:HD3	1.82	0.41
1:A:283:GLN:HG3	1:A:284:THR:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:553:ASN:C	2:B:557:GLU:HG2	2.41	0.41
2:B:635:ASN:HD22	2:B:636:GLY:N	2.19	0.41
1:A:79:SER:HB3	1:A:80:TRP:H	1.77	0.41
1:A:127:TRP:N	1:A:128:PRO:CD	2.84	0.40
1:A:5:ASP:HA	2:B:527:GLN:O	2.21	0.40
2:B:635:ASN:ND2	2:B:635:ASN:H	2.19	0.40
1:A:132:THR:HB	1:A:152:LEU:HD11	2.03	0.40
2:B:573:LEU:HA	2:B:573:LEU:HD23	1.81	0.40
1:A:120:ILE:HG23	1:A:121:PHE:N	2.36	0.40
1:A:139:CYS:O	1:A:146:SER:HB3	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:GLN:NE2	6:B:31:HOH:O[9_555]	2.18	0.02

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	322/324 (99%)	278 (86%)	31 (10%)	13 (4%)	3	17
2	B	158/160 (99%)	141 (89%)	16 (10%)	1 (1%)	28	67
All	All	480/484 (99%)	419 (87%)	47 (10%)	14 (3%)	5	25

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	80	TRP
2	B	634	GLY
1	A	27	ASN

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Mol	Chain	Res	Type
1	A	79	SER
1	A	134	GLY
1	A	199	ASP
1	A	125	SER
1	A	77	ALA
1	A	196	GLN
1	A	325	PRO
1	A	126	SER
1	A	210	ASN
1	A	326	SER
1	A	224	ARG

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	282/282 (100%)	263 (93%)	19 (7%)	19	52
2	B	136/136 (100%)	122 (90%)	14 (10%)	8	29
All	All	418/418 (100%)	385 (92%)	33 (8%)	14	43

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ILE
1	A	35	ASN
1	A	38	GLU
1	A	42	ASN
1	A	48	LEU
1	A	79	SER
1	A	91	ASN
1	A	99	PHE
1	A	126	SER
1	A	142	PHE
1	A	188	THR
1	A	193	THR

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Mol	Chain	Res	Type
1	A	214	THR
1	A	261	ASN
1	A	270	THR
1	A	271	SER
1	A	280	THR
1	A	296	GLN
1	A	321	LEU
2	B	506	MET
2	B	515	THR
2	B	538	GLN
2	B	560	ASN
2	B	562	GLN
2	B	572	ASN
2	B	576	ARG
2	B	580	LEU
2	B	590	ASP
2	B	603	GLU
2	B	616	ARG
2	B	620	GLU
2	B	635	ASN
2	B	652	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	35	ASN
1	A	42	ASN
1	A	191	GLN
1	A	192	GLN
1	A	226	GLN
1	A	250	ASN
1	A	261	ASN
1	A	276	HIS
1	A	290	ASN
1	A	299	HIS
2	B	530	GLN
2	B	553	ASN
2	B	560	ASN
2	B	572	ASN
2	B	625	GLN
2	B	628	ASN

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Mol	Chain	Res	Type
2	B	635	ASN
2	B	642	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

5 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	$\# Z > 2$
5	SIA	A	3033	5	17,20,21	0.41	0	19,28,31	0.78	1 (5%)
5	GAL	A	3034	5	11,11,12	0.48	0	13,15,17	0.73	1 (7%)
5	GLC	A	3035	5	12,12,12	0.53	0	17,17,17	0.43	0
3	NAG	A	626	1,3	14,14,15	0.51	0	15,19,21	0.95	1 (6%)
3	NDG	A	627	3	14,14,15	0.75	0	15,19,21	0.78	0

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
5	SIA	A	3033	5	-	0/14/34/38	0/1/1/1
5	GAL	A	3034	5	-	0/2/19/22	0/1/1/1
5	GLC	A	3035	5	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	626	1,3	-	0/6/23/26	0/1/1/1
3	NDG	A	627	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	626	NAG	C2-N2-C7	-2.43	119.39	122.94
5	A	3033	SIA	C6-C5-N5	-2.02	107.45	111.00
5	A	3034	GAL	C1-C2-C3	2.04	112.23	109.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	3033	SIA	4	0
3	A	626	NAG	5	0
3	A	627	NDG	2	0

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	A	631	1	14,14,15	0.59	0	15,19,21	0.87	1 (6%)
4	NAG	A	632	1	14,14,15	0.83	0	15,19,21	1.13	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
4	NAG	A	631	1	-	0/6/23/26	0/1/1/1
4	NAG	A	632	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	631	NAG	C2-N2-C7	-2.15	119.81	122.94
4	A	632	NAG	C1-O5-C5	2.93	116.20	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	632	NAG	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	324/324 (100%)	-0.09	5 (1%) 74 55	28, 52, 74, 87	0
2	B	160/160 (100%)	-0.27	1 (0%) 89 77	29, 48, 68, 87	0
All	All	484/484 (100%)	-0.15	6 (1%) 79 61	28, 51, 73, 87	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	142	PHE	3.0
2	B	660	PRO	2.9
1	A	327	ILE	2.8
1	A	141	TYR	2.3
1	A	263	GLY	2.1
1	A	40	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	A	626	14/15	0.94	0.27	1.48	66,70,72,76	0
5	SIA	A	3033	20/21	0.95	0.16	-0.24	53,60,67,68	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	GAL	A	3034	11/12	0.92	0.19	-	71,77,81,83	0
5	GLC	A	3035	12/12	0.86	0.40	-	88,96,98,100	0
3	NDG	A	627	14/15	0.87	0.35	-	76,83,87,87	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(\AA^2)	Q<0.9
4	NAG	A	632	14/15	0.72	0.40	-	86,89,90,92	0
4	NAG	A	631	14/15	0.78	0.49	-	85,89,91,92	0

6.5 Other polymers [i](#)

There are no such residues in this entry.