



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:08 am GMT

PDB ID : 1HA0
Title : HEMAGGLUTININ PRECURSOR HA0
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Deposited on : 1998-10-08
Resolution : 2.80 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

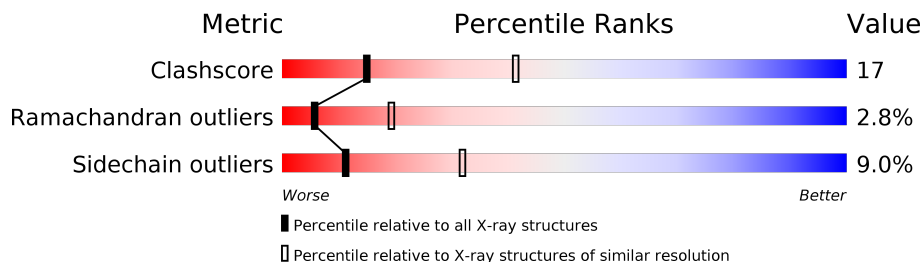
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	494	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	1440	X	-	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4070 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.

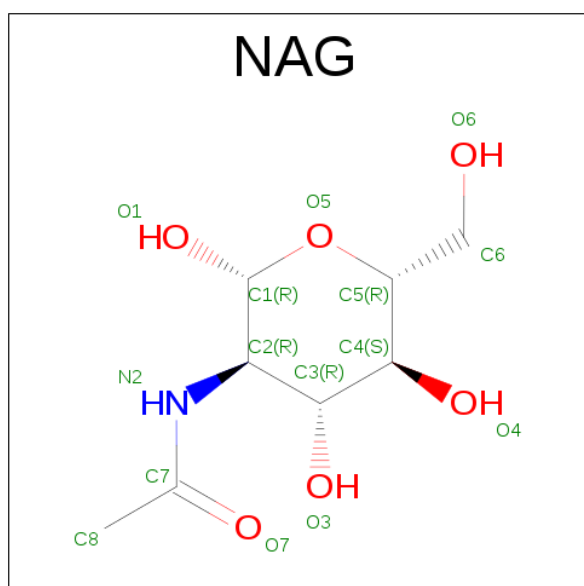
- Molecule 1 is a protein called PROTEIN (HEMAGGLUTININ PRECURSOR).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	494	3888	2426	684	759	19	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	329	GLN	ARG	ENGINEERED	UNP P03437
A	493	ASN	ASP	CONFLICT	UNP P03437

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	3	Total	C	N	O	0	0
			39	22	2	15		
3	A	3	Total	C	N	O	0	0
			39	22	2	15		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 5 is water.

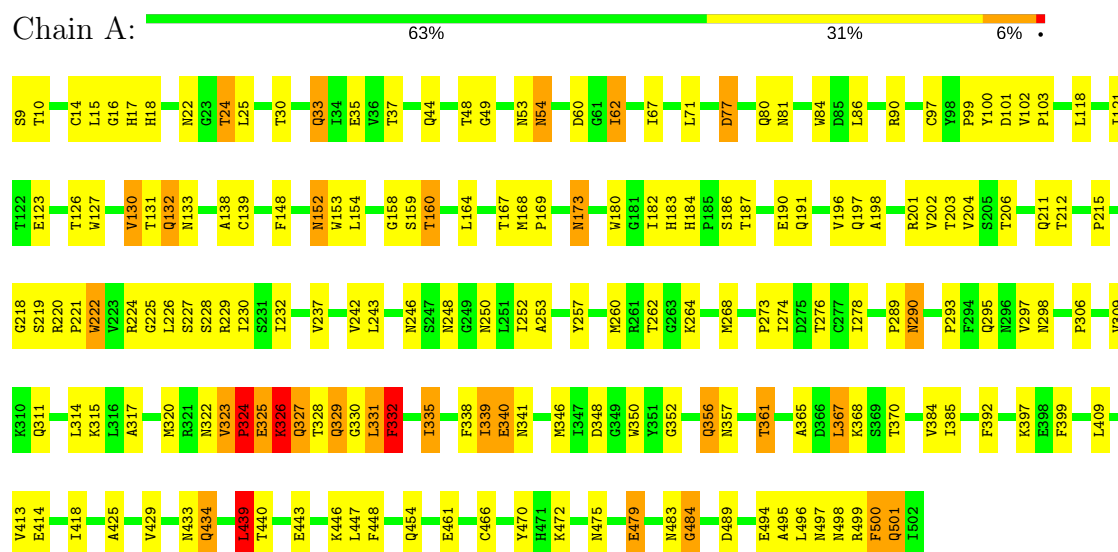
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	37	Total	O	0	0
			37	37		

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

Note EDS was not executed.

• Molecule 1: PROTEIN (HEMAGGLUTININ PRECURSOR)



4 Data and refinement statistics

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

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	153.00Å 153.00Å 153.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.80	Depositor
% Data completeness (in resolution range)	98.5 (8.00-2.80)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.224 , 0.302	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4070	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

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: BMA, NAG, MAN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.41	0/3969	0.71	4/5381 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	326	LYS	N-CA-C	6.90	129.62	111.00
1	A	327	GLN	N-CA-C	5.62	126.18	111.00
1	A	439	LEU	CA-CB-CG	5.47	127.88	115.30
1	A	328	THR	N-CA-C	5.32	125.36	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3888	0	3757	130	0
2	A	28	0	26	1	0
3	A	78	0	68	2	0
4	A	39	0	34	3	0
5	A	37	0	0	9	0
All	All	4070	0	3885	134	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 17.

All (134) 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:14:CYS:SG	1:A:335:ILE:HG22	1.86	1.14
1:A:25:LEU:HD13	1:A:33:GLN:HG3	1.47	0.96
1:A:132:GLN:HB3	1:A:152:ASN:HD21	1.30	0.94
1:A:331:LEU:O	1:A:332:PHE:HB3	1.80	0.79
1:A:173:ASN:H	1:A:173:ASN:HD22	1.31	0.77
1:A:132:GLN:HB3	1:A:152:ASN:ND2	1.99	0.76
1:A:10:THR:HG22	1:A:470:TYR:HA	1.68	0.75
1:A:182:ILE:HD11	1:A:215:PRO:HD3	1.67	0.75
1:A:335:ILE:HD12	1:A:356:GLN:HB2	1.71	0.72
1:A:293:PRO:HD3	1:A:385:ILE:HG22	1.72	0.72
1:A:187:THR:HG23	1:A:190:GLU:H	1.56	0.70
1:A:17:HIS:HE1	1:A:324:PRO:HD3	1.58	0.69
1:A:14:CYS:SG	1:A:335:ILE:CG2	2.74	0.68
1:A:500:PHE:O	1:A:501:GLN:HB2	1.93	0.68
1:A:357:ASN:HD21	1:A:475:ASN:HD22	1.44	0.66
1:A:357:ASN:HD21	1:A:475:ASN:ND2	1.94	0.66
1:A:138:ALA:HB2	1:A:226:LEU:HD11	1.79	0.65
1:A:399:PHE:HB2	5:A:1474:HOH:O	1.94	0.65
1:A:346:MET:SD	1:A:365:ALA:HB2	2.36	0.65
1:A:53:ASN:OD1	1:A:276:THR:HA	2.00	0.62
1:A:130:VAL:HG21	1:A:154:LEU:HB3	1.82	0.61
1:A:183:HIS:HB2	1:A:252:ILE:HD11	1.83	0.61
1:A:77:ASP:O	1:A:80:GLN:HG3	2.02	0.60
1:A:262:THR:HA	5:A:1486:HOH:O	2.02	0.58
1:A:325:GLU:O	1:A:326:LYS:HB2	2.03	0.58
1:A:133:ASN:H	1:A:152:ASN:HD21	1.51	0.57
1:A:203:THR:HA	1:A:211:GLN:O	2.05	0.57
1:A:290:ASN:HD22	1:A:290:ASN:H	1.51	0.57
1:A:384:VAL:HG12	1:A:385:ILE:HG23	1.88	0.56
1:A:339:ILE:HG22	1:A:340:GLU:N	2.20	0.56
1:A:201:ARG:HB3	1:A:248:ASN:OD1	2.08	0.54
1:A:127:TRP:CZ2	1:A:253:ALA:HB1	2.42	0.54
1:A:16:GLY:HA2	1:A:338:PHE:HB3	1.89	0.54
1:A:222:TRP:HA	1:A:226:LEU:O	2.08	0.53
1:A:17:HIS:CE1	1:A:324:PRO:HD3	2.41	0.53
1:A:329:GLN:HG2	1:A:330:GLY:N	2.22	0.53
1:A:121:ILE:HD12	1:A:257:TYR:OH	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:THR:HA	1:A:242:VAL:O	2.08	0.53
1:A:323:VAL:HG22	1:A:323:VAL:O	2.08	0.53
1:A:67:ILE:HG12	5:A:1463:HOH:O	2.08	0.53
1:A:317:ALA:H	1:A:433:ASN:ND2	2.07	0.52
1:A:97:CYS:O	1:A:224:ARG:NH1	2.41	0.52
1:A:320:MET:HE1	1:A:350:TRP:HB3	1.92	0.52
1:A:60:ASP:HB2	1:A:274:ILE:HD12	1.91	0.51
1:A:62:ILE:O	1:A:90:ARG:HB2	2.10	0.51
1:A:203:THR:OG1	1:A:212:THR:HG23	2.11	0.51
1:A:222:TRP:CZ3	1:A:225:GLY:HA2	2.45	0.50
1:A:309:VAL:HB	1:A:311:GLN:OE1	2.11	0.50
1:A:24:THR:HG21	1:A:315:LYS:HE2	1.93	0.49
1:A:97:CYS:HA	1:A:139:CYS:HA	1.94	0.49
1:A:290:ASN:HD22	1:A:290:ASN:N	2.06	0.48
1:A:352:GLY:HA3	1:A:365:ALA:HA	1.95	0.48
1:A:202:VAL:O	1:A:212:THR:HA	2.13	0.48
1:A:297:VAL:HG13	4:A:1460:NAG:H82	1.94	0.48
1:A:103:PRO:HD2	1:A:232:ILE:O	2.14	0.48
1:A:446:LYS:HD2	5:A:1467:HOH:O	2.14	0.48
1:A:182:ILE:HD11	1:A:215:PRO:CD	2.39	0.48
1:A:186:SER:HA	1:A:218:GLY:O	2.13	0.48
1:A:152:ASN:C	1:A:152:ASN:HD22	2.16	0.48
1:A:184:HIS:HB3	1:A:220:ARG:NH2	2.29	0.48
1:A:357:ASN:ND2	1:A:475:ASN:HD22	2.12	0.48
1:A:14:CYS:HA	1:A:466:CYS:HA	1.96	0.47
1:A:317:ALA:H	1:A:433:ASN:HD21	1.61	0.47
4:A:1461:NAG:O7	4:A:1461:NAG:C1	2.63	0.47
1:A:220:ARG:HB3	1:A:221:PRO:CD	2.45	0.47
1:A:500:PHE:O	1:A:501:GLN:CB	2.63	0.47
1:A:356:GLN:HG2	1:A:361:THR:HB	1.95	0.47
1:A:399:PHE:CD1	1:A:399:PHE:N	2.83	0.47
1:A:152:ASN:HD22	1:A:153:TRP:N	2.13	0.46
1:A:494:GLU:HG2	1:A:498:ASN:ND2	2.31	0.46
1:A:184:HIS:CE1	5:A:1487:HOH:O	2.67	0.46
1:A:290:ASN:ND2	1:A:290:ASN:H	2.13	0.46
1:A:338:PHE:HE1	1:A:448:PHE:CD1	2.34	0.46
1:A:221:PRO:O	1:A:229:ARG:NH2	2.49	0.46
1:A:260:MET:N	5:A:1480:HOH:O	2.45	0.46
1:A:331:LEU:O	1:A:332:PHE:CB	2.57	0.45
1:A:49:GLY:O	1:A:273:PRO:HD2	2.16	0.45
1:A:81:ASN:ND2	2:A:1440:NAG:C7	2.79	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:LEU:HB3	1:A:429:VAL:HG21	1.99	0.45
1:A:439:LEU:C	1:A:439:LEU:HD23	2.37	0.45
1:A:131:THR:HA	5:A:1495:HOH:O	2.15	0.45
1:A:461:GLU:HG3	5:A:1476:HOH:O	2.17	0.45
1:A:367:LEU:HA	1:A:370:THR:HB	1.99	0.45
1:A:17:HIS:HA	1:A:350:TRP:O	2.16	0.44
1:A:295:GLN:HB3	1:A:306:PRO:HB2	2.00	0.44
1:A:86:LEU:HD11	1:A:268:MET:HB2	1.99	0.44
1:A:24:THR:O	1:A:35:GLU:HA	2.17	0.44
1:A:448:PHE:HZ	1:A:461:GLU:HG3	1.82	0.44
3:A:1450:NAG:H62	3:A:1451:NAG:O5	2.17	0.44
1:A:167:THR:HA	1:A:243:LEU:O	2.17	0.44
1:A:173:ASN:H	1:A:173:ASN:ND2	2.08	0.43
1:A:237:VAL:HG21	1:A:243:LEU:HB2	2.00	0.43
1:A:479:GLU:O	1:A:483:ASN:ND2	2.51	0.43
1:A:102:VAL:HG22	1:A:232:ILE:HB	2.00	0.43
1:A:197:GLN:HG3	1:A:248:ASN:O	2.18	0.43
1:A:44:GLN:OE1	1:A:289:PRO:HG2	2.19	0.43
1:A:324:PRO:O	1:A:325:GLU:HB2	2.18	0.43
1:A:15:LEU:CD1	1:A:447:LEU:HG	2.48	0.43
1:A:220:ARG:HD2	1:A:220:ARG:N	2.34	0.43
1:A:169:PRO:HA	1:A:242:VAL:HG23	2.01	0.43
1:A:126:THR:HG23	1:A:126:THR:O	2.19	0.43
1:A:10:THR:CG2	1:A:470:TYR:HA	2.43	0.43
1:A:298:ASN:HA	5:A:1481:HOH:O	2.18	0.43
1:A:409:LEU:O	1:A:413:VAL:HG23	2.19	0.43
1:A:123:GLU:HB3	1:A:168:MET:HE3	2.00	0.43
1:A:84:TRP:CZ3	1:A:118:LEU:HG	2.54	0.42
1:A:37:THR:HG22	1:A:322:ASN:OD1	2.19	0.42
1:A:54:ASN:O	1:A:278:ILE:HA	2.19	0.42
1:A:434:GLN:HA	1:A:434:GLN:HE21	1.85	0.42
1:A:494:GLU:HG2	1:A:498:ASN:HD22	1.84	0.42
1:A:218:GLY:O	1:A:220:ARG:NH1	2.53	0.42
1:A:123:GLU:HB3	1:A:168:MET:CE	2.49	0.42
1:A:164:LEU:O	1:A:246:ASN:HA	2.20	0.42
1:A:454:GLN:NE2	1:A:484:GLY:HA2	2.34	0.41
1:A:160:THR:HA	1:A:196:VAL:HG21	2.02	0.41
1:A:414:GLU:O	1:A:418:ILE:HG13	2.19	0.41
3:A:1450:NAG:H3	3:A:1450:NAG:O7	2.20	0.41
1:A:100:TYR:HB2	1:A:230:ILE:O	2.21	0.41
1:A:173:ASN:HD22	1:A:173:ASN:N	2.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:495:ALA:O	1:A:499:ARG:HG3	2.20	0.41
1:A:180:TRP:CZ2	1:A:204:VAL:HB	2.55	0.41
1:A:220:ARG:HD3	1:A:227:SER:O	2.21	0.41
1:A:99:PRO:O	1:A:229:ARG:HD3	2.21	0.41
1:A:35:GLU:OE1	1:A:326:LYS:HD3	2.21	0.41
1:A:84:TRP:HZ3	1:A:118:LEU:HG	1.86	0.41
1:A:191:GLN:HG2	1:A:198:ALA:O	2.20	0.41
1:A:443:GLU:HA	1:A:446:LYS:HE3	2.01	0.41
4:A:1461:NAG:H62	4:A:1462:MAN:C1	2.51	0.41
1:A:138:ALA:CB	1:A:226:LEU:HD11	2.49	0.40
1:A:314:LEU:CD1	1:A:425:ALA:HB1	2.51	0.40
1:A:71:LEU:O	1:A:148:PHE:HB3	2.21	0.40
1:A:320:MET:HB3	1:A:440:THR:HB	2.03	0.40
1:A:215:PRO:HG3	1:A:250:ASN:ND2	2.36	0.40
1:A:264:LYS:O	1:A:392:PHE:HB3	2.22	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	492/494 (100%)	426 (87%)	52 (11%)	14 (3%)	6 19

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	GLN
1	A	324	PRO
1	A	327	GLN
1	A	331	LEU
1	A	332	PHE
1	A	339	ILE

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Mol	Chain	Res	Type
1	A	501	GLN
1	A	326	LYS
1	A	62	ILE
1	A	484	GLY
1	A	500	PHE
1	A	335	ILE
1	A	130	VAL
1	A	158	GLY

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	431/431 (100%)	392 (91%)	39 (9%)	11	32

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	18	HIS
1	A	22	ASN
1	A	24	THR
1	A	30	THR
1	A	33	GLN
1	A	48	THR
1	A	54	ASN
1	A	77	ASP
1	A	101	ASP
1	A	152	ASN
1	A	159	SER
1	A	160	THR
1	A	173	ASN
1	A	219	SER
1	A	222	TRP
1	A	228	SER
1	A	290	ASN

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Mol	Chain	Res	Type
1	A	323	VAL
1	A	324	PRO
1	A	325	GLU
1	A	326	LYS
1	A	329	GLN
1	A	332	PHE
1	A	340	GLU
1	A	341	ASN
1	A	348	ASP
1	A	356	GLN
1	A	361	THR
1	A	367	LEU
1	A	368	LYS
1	A	397	LYS
1	A	434	GLN
1	A	439	LEU
1	A	472	LYS
1	A	479	GLU
1	A	489	ASP
1	A	496	LEU
1	A	497	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	33	GLN
1	A	54	ASN
1	A	132	GLN
1	A	152	ASN
1	A	173	ASN
1	A	191	GLN
1	A	211	GLN
1	A	246	ASN
1	A	250	ASN
1	A	290	ASN
1	A	424	ASN
1	A	433	ASN
1	A	434	GLN
1	A	454	GLN
1	A	475	ASN
1	A	483	ASN

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Mol	Chain	Res	Type
1	A	498	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 ⓘ

9 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	1430	1,3	14,14,15	0.65	0	15,19,21	1.71	3 (20%)
3	NAG	A	1431	3	14,14,15	0.44	0	15,19,21	0.78	0
3	BMA	A	1432	3	11,11,12	0.42	0	13,15,17	0.51	0
3	NAG	A	1450	1,3	14,14,15	0.86	1 (7%)	15,19,21	1.85	2 (13%)
3	NAG	A	1451	3	14,14,15	0.73	0	15,19,21	1.66	2 (13%)
3	BMA	A	1452	3	11,11,12	0.48	0	13,15,17	0.43	0
4	NAG	A	1460	1,4	14,14,15	0.79	1 (7%)	15,19,21	0.72	0
4	NAG	A	1461	4	14,14,15	0.73	0	15,19,21	1.53	3 (20%)
4	MAN	A	1462	4	11,11,12	0.74	0	13,15,17	1.04	1 (7%)

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	1430	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1431	3	-	0/6/23/26	0/1/1/1
3	BMA	A	1432	3	-	0/2/19/22	0/1/1/1
3	NAG	A	1450	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1451	3	-	0/6/23/26	0/1/1/1
3	BMA	A	1452	3	-	0/2/19/22	0/1/1/1
4	NAG	A	1460	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	1461	4	-	0/6/23/26	0/1/1/1
4	MAN	A	1462	4	-	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	A	1450	NAG	C1-C2	2.04	1.55	1.52
4	A	1460	NAG	C1-C2	2.51	1.55	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1450	NAG	C4-C3-C2	-5.72	102.64	111.02
3	A	1451	NAG	C4-C3-C2	-5.05	103.61	111.02
3	A	1430	NAG	C4-C3-C2	-4.75	104.06	111.02
4	A	1461	NAG	C4-C3-C2	-3.77	105.49	111.02
4	A	1461	NAG	O5-C1-C2	-2.70	107.71	111.47
3	A	1430	NAG	O5-C1-C2	-2.67	107.75	111.47
3	A	1451	NAG	O5-C1-C2	-2.38	108.17	111.47
3	A	1430	NAG	C1-C2-N2	2.37	114.53	110.49
3	A	1450	NAG	C1-O5-C5	2.44	115.54	112.17
4	A	1461	NAG	C1-C2-N2	2.84	115.35	110.49
4	A	1462	MAN	C1-O5-C5	2.95	116.24	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1450	NAG	2	0
3	A	1451	NAG	1	0
4	A	1460	NAG	1	0
4	A	1461	NAG	2	0
4	A	1462	MAN	1	0

5.6 Ligand geometry

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
2	NAG	A	1420	1	14,14,15	0.86	1 (7%)	15,19,21	1.66	3 (20%)
2	NAG	A	1440	1	14,14,15	0.50	0	15,19,21	1.74	2 (13%)

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	NAG	A	1420	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1440	1	1/1/5/7	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1420	NAG	O5-C5	2.13	1.47	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1440	NAG	C2-N2-C7	-2.12	119.85	122.94
2	A	1420	NAG	C4-C3-C2	-2.03	108.04	111.02
2	A	1420	NAG	O5-C1-C2	2.40	114.81	111.47
2	A	1420	NAG	C1-O5-C5	4.60	118.50	112.17
2	A	1440	NAG	O5-C1-C2	5.71	119.42	111.47

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1440	NAG	C1

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1440	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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.