



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 08:30 am GMT

PDB ID : 1E8T  
Title : STRUCTURE OF THE MULTIFUNCTIONAL PARAMYXOVIRUS HEMA  
GGLUTININ-NEURAMINIDASE  
Authors : Crennell, S.; Takimoto, T.; Portner, A.; Taylor, G.  
Deposited on : 2000-10-01  
Resolution : 2.50 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](mailto: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.

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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

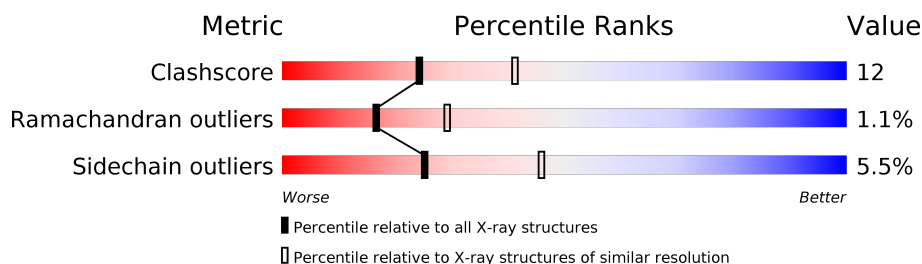
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)

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  $\geq 3$ , 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	454	
1	B	454	

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	GOL	A	1003	-	-	X	-

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7214 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.

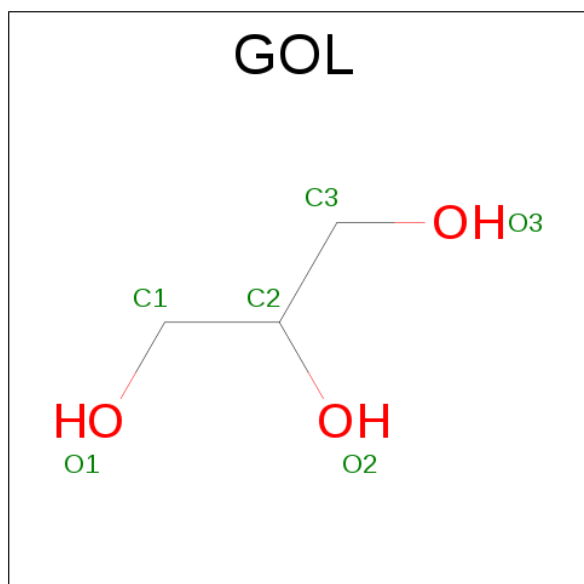
- Molecule 1 is a protein called HEMAGGLUTININ-NEURAMINIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	447	Total	C	N	O	S	0	0	1
			3446	2172	588	667	19			
1	B	450	Total	C	N	O	S	0	0	1
			3468	2185	594	670	19			

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

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



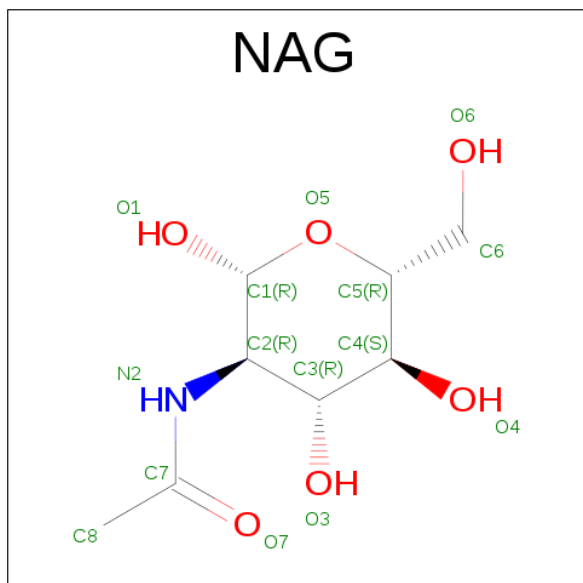
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

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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		

- 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		

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

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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	140	Total	O	0	0
			140	140		

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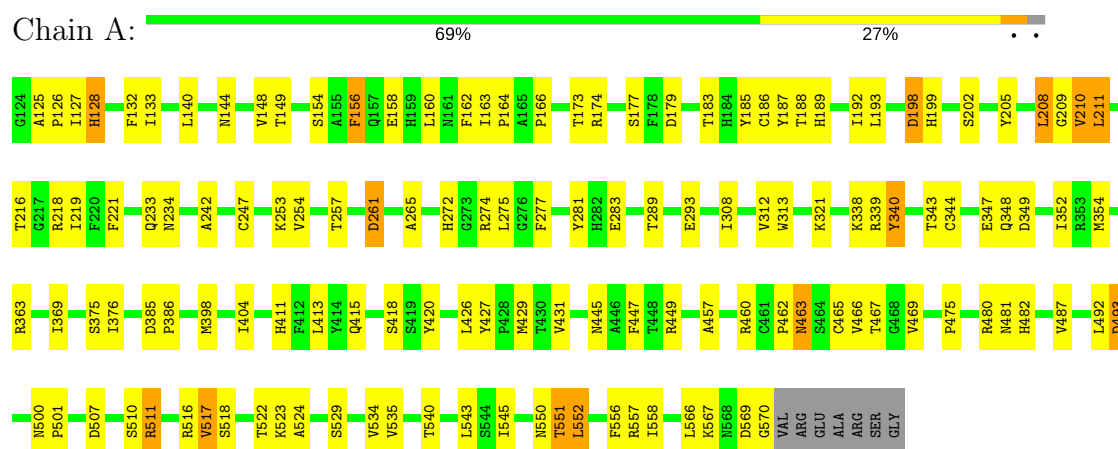
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	71	Total	O	0	0
			71	71		

### 3 Residue-property plots

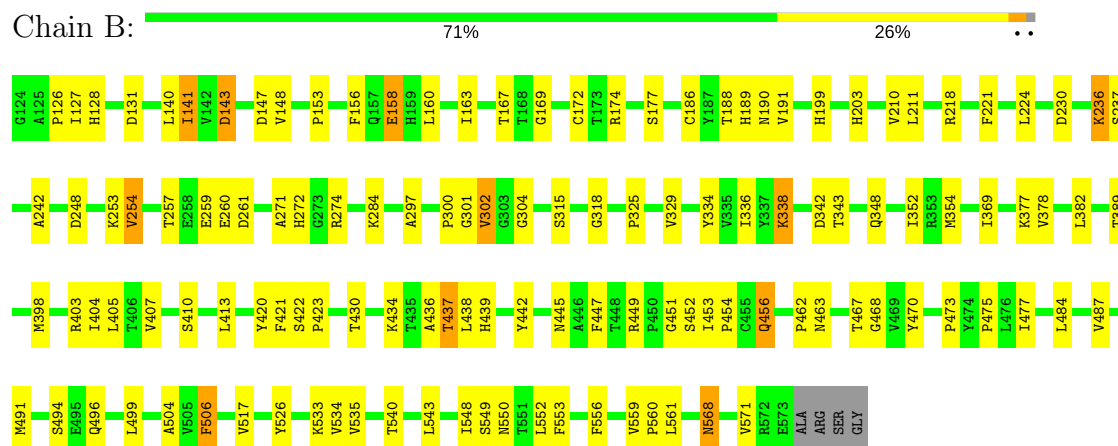
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

#### • Molecule 1: HEMAGGLUTININ-NEURAMINIDASE



#### • Molecule 1: HEMAGGLUTININ-NEURAMINIDASE



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.71Å 77.91Å 198.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.50	Depositor
% Data completeness (in resolution range)	97.0 (6.00-2.50)	Depositor
$R_{merge}$	0.03	Depositor
$R_{sym}$	0.03	Depositor
Refinement program	CNS	Depositor
R, $R_{free}$	0.222 , 0.276	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7214	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	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: GOL, CA, NAG

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	0.40	0/3531	0.70	0/4804
1	B	0.35	0/3553	0.62	0/4833
All	All	0.38	0/7084	0.66	0/9637

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	3446	0	3340	92	0
1	B	3468	0	3366	80	0
2	A	1	0	0	0	0
3	A	12	0	16	5	0
3	B	6	0	8	1	0
4	A	14	0	13	2	0
5	A	28	0	25	0	0
5	B	28	0	25	0	0
6	A	140	0	0	3	0
6	B	71	0	0	4	0
All	All	7214	0	6793	171	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 12.

All (171) 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:511:ARG:HH11	1:A:511:ARG:HG2	1.30	0.96
1:B:156:PHE:HE2	1:B:517:VAL:HA	1.35	0.91
1:A:128:HIS:HA	1:A:210:VAL:HG13	1.54	0.87
1:A:128:HIS:CD2	1:A:211:LEU:H	1.96	0.83
1:B:156:PHE:CE2	1:B:517:VAL:HA	2.18	0.78
1:A:156:PHE:CE2	1:A:517:VAL:HA	2.23	0.73
1:A:404:ILE:HD13	1:A:413:LEU:HD23	1.71	0.73
1:B:284:LYS:HD3	1:B:382:LEU:O	1.90	0.72
1:B:128:HIS:CD2	1:B:211:LEU:H	2.08	0.71
1:B:128:HIS:HD2	1:B:211:LEU:H	1.40	0.70
1:B:549:SER:HB3	1:B:556:PHE:HA	1.74	0.69
1:A:511:ARG:HH11	1:A:511:ARG:CG	2.05	0.68
1:A:149:THR:HG21	1:A:510:SER:HB2	1.76	0.67
1:B:421:PHE:CE2	1:B:423:PRO:HG2	2.29	0.66
1:B:177:SER:HB3	1:B:188:THR:HG22	1.80	0.64
1:A:469:VAL:HG11	1:A:492:LEU:HD22	1.80	0.64
1:A:466:VAL:H	3:A:1003:GOL:H11	1.62	0.64
1:A:192:ILE:HD13	1:A:202:SER:HB3	1.80	0.62
1:A:462:PRO:HD2	6:A:2052:HOH:O	1.99	0.62
1:A:154:SER:HB3	1:A:566:LEU:HD11	1.82	0.60
1:A:501:PRO:HG3	1:A:524:ALA:HB2	1.84	0.60
1:A:218:ARG:NE	1:A:218:ARG:HA	2.16	0.59
1:A:339:ARG:NH2	1:A:457:ALA:O	2.35	0.59
1:B:257:THR:HB	1:B:260:GLU:HB3	1.83	0.59
1:B:221:PHE:CE2	1:B:543:LEU:HD22	2.38	0.59
1:B:430:THR:OG1	1:B:437:THR:HG23	2.03	0.58
1:A:338:LYS:HZ1	1:A:344:CYS:HB3	1.67	0.58
1:A:369:ILE:HB	1:A:398:MET:HG3	1.86	0.57
1:A:340:TYR:HE2	4:A:1570:NAG:H82	1.70	0.57
1:A:177:SER:HB3	1:A:188:THR:HG22	1.86	0.56
1:A:179:ASP:HA	6:A:2013:HOH:O	2.04	0.56
1:A:156:PHE:HE2	1:A:517:VAL:HA	1.69	0.55
1:B:297:ALA:O	1:B:318:GLY:HA3	2.07	0.54
1:A:465:CYS:HA	3:A:1003:GOL:H32	1.91	0.53
1:A:156:PHE:CZ	1:A:517:VAL:HA	2.44	0.53
1:B:421:PHE:CZ	1:B:423:PRO:HG2	2.43	0.53
1:B:237:SER:O	1:B:300:PRO:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:ASN:HD21	1:A:516:ARG:NH1	2.06	0.52
1:B:334:TYR:OH	1:B:352:ILE:HD12	2.10	0.52
1:B:420:TYR:CG	1:B:462:PRO:HA	2.45	0.52
1:A:265:ALA:O	1:A:321:LYS:HE2	2.09	0.52
1:B:302:VAL:HG11	1:B:526:TYR:CE2	2.45	0.52
1:A:218:ARG:HE	1:A:219:ILE:N	2.09	0.51
1:A:308:ILE:HG21	1:A:431:VAL:HG21	1.91	0.51
1:A:218:ARG:HE	1:A:219:ILE:H	1.58	0.51
1:A:511:ARG:NH1	1:A:511:ARG:HG2	2.08	0.51
1:A:339:ARG:HD2	1:A:445:ASN:OD1	2.11	0.50
1:B:451:GLY:O	1:B:456:GLN:HA	2.11	0.50
1:B:405:LEU:H	1:B:405:LEU:HD23	1.77	0.50
1:B:534:VAL:O	1:B:534:VAL:HG23	2.13	0.49
1:B:533:LYS:HG2	1:B:534:VAL:N	2.28	0.49
1:A:340:TYR:CE2	4:A:1570:NAG:H82	2.48	0.49
1:B:491:MET:CE	1:B:506:PHE:HE2	2.25	0.49
1:A:234:ASN:HB2	1:A:254:VAL:O	2.13	0.48
1:A:466:VAL:N	3:A:1003:GOL:H11	2.28	0.48
1:A:140:LEU:HG	1:A:540:THR:HG23	1.96	0.48
1:B:191:VAL:HG13	1:B:203:HIS:HB2	1.96	0.48
1:B:494:SER:HB2	1:B:499:LEU:HD12	1.94	0.48
1:A:132:PHE:CZ	1:A:211:LEU:HD22	2.49	0.48
1:A:518:SER:HB3	1:A:522:THR:OG1	2.13	0.48
1:A:173:THR:HB	1:A:558:ILE:HD12	1.96	0.48
1:B:405:LEU:HD23	1:B:405:LEU:N	2.29	0.48
1:A:133:ILE:O	1:A:535:VAL:HG21	2.14	0.48
1:A:160:LEU:HD13	1:A:221:PHE:CD1	2.49	0.48
1:B:548:ILE:HD11	1:B:559:VAL:HG21	1.95	0.47
1:B:148:VAL:HG13	1:B:477:ILE:HD12	1.96	0.47
1:A:411:HIS:HB2	1:A:429:MET:O	2.15	0.47
1:B:491:MET:HE2	1:B:506:PHE:CE2	2.49	0.47
1:A:185:TYR:CZ	1:A:209:GLY:HA3	2.50	0.47
1:A:148:VAL:HG11	1:A:507:ASP:HB3	1.97	0.46
1:A:385:ASP:N	1:A:386:PRO:HD3	2.31	0.46
1:A:551:THR:OG1	1:A:552:LEU:HD22	2.14	0.46
1:B:140:LEU:HD13	1:B:540:THR:HG23	1.98	0.46
1:A:281:TYR:CE2	1:A:283:GLU:HB2	2.51	0.46
1:B:236:LYS:HB2	1:B:236:LYS:NZ	2.31	0.46
1:B:274:ARG:HH21	1:B:378:VAL:HG13	1.79	0.46
1:B:163:ILE:HD12	1:B:560:PRO:HB3	1.96	0.46
1:B:467:THR:HG23	1:B:468:GLY:N	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:568:ASN:HB2	1:B:571:VAL:HG13	1.96	0.46
1:A:219:ILE:HD11	1:A:534:VAL:HG21	1.98	0.46
1:A:363:ARG:H	3:A:1003:GOL:HO2	1.62	0.46
1:A:354:MET:HE2	1:A:463:ASN:HA	1.98	0.46
1:B:453:ILE:HG13	1:B:496:GLN:NE2	2.30	0.46
1:A:289:THR:O	1:A:293:GLU:HB3	2.16	0.46
1:A:493:ASP:HB3	1:A:500:ASN:ND2	2.31	0.46
1:B:517:VAL:HG13	1:B:561:LEU:HD11	1.96	0.45
1:A:481:ASN:O	1:A:482:HIS:HB2	2.16	0.45
1:B:224:LEU:HD23	6:B:2020:HOH:O	2.16	0.45
1:A:511:ARG:NH1	1:A:511:ARG:CG	2.75	0.45
1:B:158:GLU:HG2	1:B:218:ARG:HG3	1.99	0.45
1:A:415:GLN:NE2	1:A:427:TYR:OH	2.49	0.45
1:A:308:ILE:CG2	1:A:431:VAL:HG21	2.47	0.45
1:B:452:SER:HB2	6:B:2061:HOH:O	2.17	0.45
1:A:186:CYS:SG	1:A:242:ALA:HA	2.56	0.45
1:A:166:PRO:O	1:B:169:GLY:HA3	2.17	0.45
1:A:475:PRO:HA	1:A:487:VAL:HG12	1.99	0.45
1:B:550:ASN:HB3	1:B:553:PHE:HD2	1.82	0.44
1:A:543:LEU:HD21	1:A:545:ILE:HD11	2.00	0.44
1:B:423:PRO:HB3	1:B:445:ASN:HA	1.98	0.44
1:B:272:HIS:CE1	1:B:382:LEU:HD23	2.52	0.44
1:B:325:PRO:O	1:B:329:VAL:HG23	2.17	0.44
1:B:403:ARG:HB2	1:B:470:TYR:OH	2.18	0.44
1:B:413:LEU:HD13	1:B:438:LEU:HD11	1.99	0.44
1:A:257:THR:O	1:A:261:ASP:HB2	2.17	0.44
1:B:174:ARG:NH2	1:B:190:ASN:OD1	2.48	0.44
1:B:211:LEU:HD21	1:B:221:PHE:CE1	2.53	0.44
1:B:475:PRO:HA	1:B:487:VAL:HG12	1.99	0.44
1:B:127:ILE:HG12	1:B:128:HIS:N	2.33	0.44
1:B:389:THR:HG21	1:B:436:ALA:O	2.18	0.43
1:A:420:TYR:CG	1:A:462:PRO:HA	2.53	0.43
1:A:233:GLN:NE2	1:A:253:LYS:HZ3	2.17	0.43
1:B:439:HIS:ND1	3:B:1004:GOL:H32	2.33	0.43
1:B:141:ILE:HD11	1:B:143:ASP:HB2	2.00	0.43
1:B:369:ILE:HB	1:B:398:MET:HG3	1.99	0.43
1:B:422:SER:N	1:B:423:PRO:HD2	2.33	0.43
1:A:187:TYR:OH	1:A:558:ILE:HG21	2.19	0.43
1:B:407:VAL:HG21	1:B:475:PRO:HB2	1.99	0.43
1:A:133:ILE:O	1:A:133:ILE:HG13	2.19	0.42
1:A:208:LEU:HD22	1:A:247:CYS:SG	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:HIS:HD2	1:A:211:LEU:H	1.58	0.42
1:A:274:ARG:O	1:A:281:TYR:HA	2.19	0.42
1:A:154:SER:HB3	1:A:566:LEU:CD1	2.49	0.42
1:A:529:SER:HA	1:A:543:LEU:O	2.19	0.42
1:A:144:ASN:HD21	1:A:480:ARG:H	1.67	0.42
1:A:348:GLN:O	1:A:352:ILE:HG13	2.19	0.42
1:B:172:CYS:HB2	1:B:556:PHE:O	2.19	0.42
1:A:185:TYR:CE2	1:A:221:PHE:HD2	2.37	0.42
1:B:301:GLY:O	1:B:302:VAL:HB	2.20	0.42
1:A:567:LYS:C	1:A:569:ASP:H	2.23	0.42
1:A:308:ILE:HB	1:A:313:TRP:NE1	2.35	0.42
1:A:523:LYS:HG3	6:A:2114:HOH:O	2.20	0.42
1:A:426:LEU:HA	1:A:426:LEU:HD23	1.88	0.42
1:B:126:PRO:HD3	6:B:2001:HOH:O	2.18	0.42
1:B:254:VAL:HG11	1:B:261:ASP:OD2	2.20	0.42
1:B:315:SER:HB3	1:B:404:ILE:HG13	2.02	0.42
1:B:354:MET:HE2	1:B:463:ASN:HA	2.02	0.42
1:B:160:LEU:HD13	1:B:221:PHE:CD1	2.55	0.41
1:A:125:ALA:HB1	1:A:183:THR:HG21	2.01	0.41
1:B:442:TYR:CZ	1:B:484:LEU:HB3	2.55	0.41
1:A:552:LEU:H	1:A:552:LEU:HD22	1.85	0.41
1:B:131:ASP:O	1:B:535:VAL:HG22	2.19	0.41
1:B:153:PRO:HG2	1:B:504:ALA:HA	2.02	0.41
1:B:491:MET:HE2	1:B:506:PHE:HE2	1.83	0.41
1:B:403:ARG:HA	6:B:2035:HOH:O	2.20	0.41
1:A:312:VAL:O	1:A:375:SER:HA	2.20	0.41
1:A:523:LYS:HB3	1:A:556:PHE:CE1	2.56	0.41
1:B:186:CYS:SG	1:B:242:ALA:HA	2.60	0.41
1:B:203:HIS:ND1	1:B:230:ASP:HA	2.35	0.41
1:A:418:SER:HA	1:A:467:THR:O	2.20	0.41
1:B:304:GLY:N	1:B:403:ARG:HG3	2.36	0.41
1:A:148:VAL:CG1	1:A:507:ASP:HB3	2.50	0.41
1:B:336:ILE:CD1	1:B:352:ILE:HD13	2.51	0.41
1:A:174:ARG:HH12	1:A:198:ASP:HA	1.86	0.41
1:B:348:GLN:O	1:B:352:ILE:HG12	2.21	0.41
1:B:377:LYS:HA	1:B:377:LYS:HD2	1.94	0.41
1:B:210:VAL:HG13	1:B:224:LEU:HD11	2.02	0.41
1:B:253:LYS:HE3	1:B:271:ALA:HB2	2.03	0.41
1:B:506:PHE:CD1	1:B:506:PHE:N	2.88	0.41
1:A:162:PHE:CD1	1:A:221:PHE:HB2	2.56	0.41
1:A:163:ILE:HA	1:A:164:PRO:HD3	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:404:ILE:CD1	1:A:413:LEU:HD23	2.48	0.41
1:A:164:PRO:HG2	1:A:205:TYR:CZ	2.57	0.40
1:A:166:PRO:HD3	1:A:557:ARG:HG3	2.03	0.40
1:A:126:PRO:HD3	1:A:277:PHE:CE1	2.56	0.40
1:A:344:CYS:HB2	1:A:460:ARG:O	2.21	0.40
1:A:465:CYS:HA	3:A:1003:GOL:C3	2.52	0.40
1:B:257:THR:HB	1:B:260:GLU:CB	2.49	0.40
1:A:272:HIS:ND1	1:A:376:ILE:HG21	2.36	0.40
1:B:334:TYR:CZ	1:B:352:ILE:HG23	2.56	0.40
1:A:567:LYS:HE3	1:A:570:GLY:N	2.37	0.40
1:B:338:LYS:H	1:B:338:LYS:CD	2.34	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	445/454 (98%)	406 (91%)	33 (7%)	6 (1%)	14	25
1	B	448/454 (99%)	410 (92%)	34 (8%)	4 (1%)	20	36
All	All	893/908 (98%)	816 (91%)	67 (8%)	10 (1%)	17	29

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	128	HIS
1	A	551	THR
1	B	158	GLU
1	A	127	ILE
1	A	156	PHE
1	A	340	TYR
1	A	493	ASP

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Mol	Chain	Res	Type
1	B	342	ASP
1	B	302	VAL
1	B	473	PRO

### 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	387/392 (99%)	366 (95%)	21 (5%)	26	47
1	B	389/392 (99%)	367 (94%)	22 (6%)	24	44
All	All	776/784 (99%)	733 (94%)	43 (6%)	25	46

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	158	GLU
1	A	189	HIS
1	A	193	LEU
1	A	198	ASP
1	A	199	HIS
1	A	208	LEU
1	A	210	VAL
1	A	211	LEU
1	A	216	THR
1	A	261	ASP
1	A	275	LEU
1	A	343	THR
1	A	347	GLU
1	A	349	ASP
1	A	447	PHE
1	A	449	ARG
1	A	463	ASN
1	A	511	ARG
1	A	517	VAL
1	A	550	ASN

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Mol	Chain	Res	Type
1	A	552	LEU
1	B	141	ILE
1	B	143	ASP
1	B	147	ASP
1	B	167	THR
1	B	189	HIS
1	B	199	HIS
1	B	236	LYS
1	B	248	ASP
1	B	254	VAL
1	B	259	GLU
1	B	338	LYS
1	B	343	THR
1	B	410	SER
1	B	434	LYS
1	B	437	THR
1	B	447	PHE
1	B	449	ARG
1	B	454	PRO
1	B	456	GLN
1	B	506	PHE
1	B	552	LEU
1	B	568	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	HIS
1	A	199	HIS
1	A	203	HIS
1	A	233	GLN
1	A	323	ASN
1	A	371	GLN
1	A	415	GLN
1	A	463	ASN
1	A	482	HIS
1	A	500	ASN
1	B	128	HIS
1	B	157	GLN
1	B	351	GLN
1	B	496	GLN

### 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 ⓘ

4 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	NAG	A	1571	1,5	14,14,15	0.43	0	15,19,21	0.72	0
5	NAG	A	1572	5	14,14,15	0.43	0	15,19,21	0.61	0
5	NAG	B	1573	1,5	14,14,15	0.42	0	15,19,21	0.68	0
5	NAG	B	1574	5	14,14,15	0.40	0	15,19,21	0.74	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	NAG	A	1571	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	1572	5	-	0/6/23/26	0/1/1/1
5	NAG	B	1573	1,5	-	0/6/23/26	0/1/1/1
5	NAG	B	1574	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.



There are no ring outliers.

No monomer is involved in short contacts.

## 5.6 Ligand geometry

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	1002	-	5,5,5	0.52	0	5,5,5	0.34	0
3	GOL	A	1003	-	5,5,5	1.18	1 (20%)	5,5,5	0.34	0
4	NAG	A	1570	1	14,14,15	0.41	0	15,19,21	0.75	0
3	GOL	B	1004	-	5,5,5	0.95	0	5,5,5	0.23	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
3	GOL	A	1002	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1003	-	-	0/4/4/4	0/0/0/0
4	NAG	A	1570	1	-	0/6/23/26	0/1/1/1
3	GOL	B	1004	-	-	0/4/4/4	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1003	GOL	C1-C2	2.19	1.60	1.52

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1003	GOL	5	0
4	A	1570	NAG	2	0
3	B	1004	GOL	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.