



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:10 PM GMT

PDB ID : 4HF5
Title : Crystal structure of Fab 8F8 in complex a H2N2 influenza virus hemagglutinin
Authors : Xu, R.; Wilson, I.A.
Deposited on : 2012-10-04
Resolution : 3.00 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

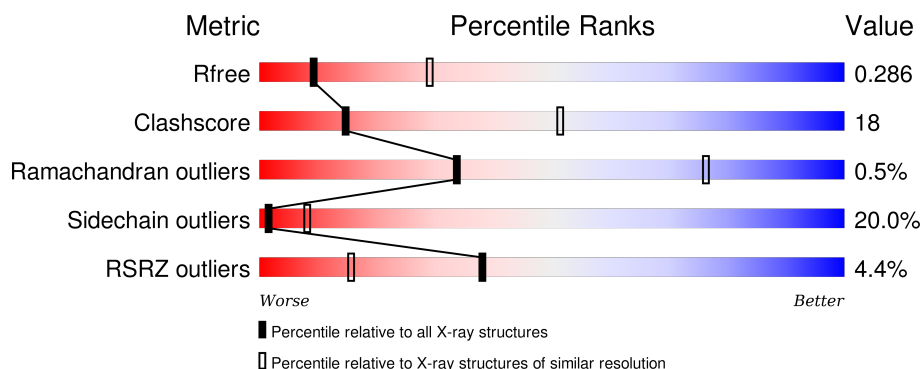
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	327	<div> <div>64%</div> <div>29%</div> <div>6%</div> <div>.</div> </div>
2	B	174	<div> <div>%</div> <div>57%</div> <div>32%</div> <div>9%</div> <div>..</div> </div>
3	H	233	<div> <div>3%</div> <div>34%</div> <div>27%</div> <div>.</div> <div>34%</div> </div>
4	L	218	<div> <div>12%</div> <div>40%</div> <div>38%</div> <div>11%</div> <div>.</div> <div>10%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6639 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 HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2534	1593	440	486	15			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	PRO	-	EXPRESSION TAG	UNP C7S226

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	172	Total	C	N	O	S	0	0	0
			1396	871	237	279	9			

- Molecule 3 is a protein called Fab 8F8 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	153	Total	C	N	O	S	0	0	0
			1195	756	203	228	8			

- Molecule 4 is a protein called Fab 8F8 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	196	Total	C	N	O	S	0	0	0
			1436	897	237	298	4			

- Molecule 5 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	4	Total	C	N	O	0	0
			50	28	2	20		

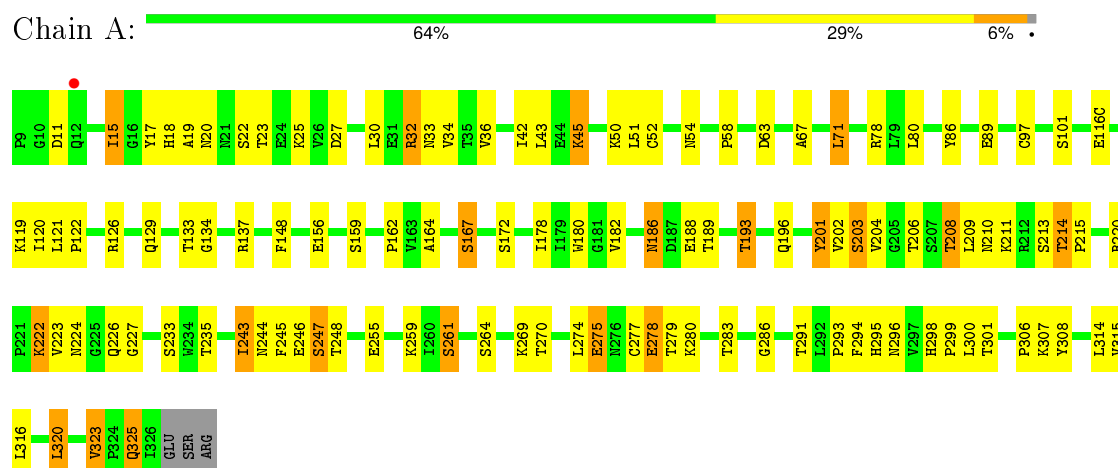
- Molecule 6 is a polymer of unknown type called SUGAR (2-MER).

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

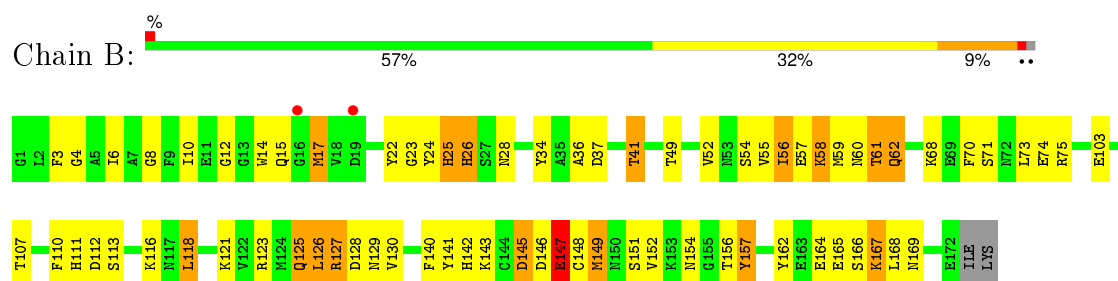
3 Residue-property plots

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

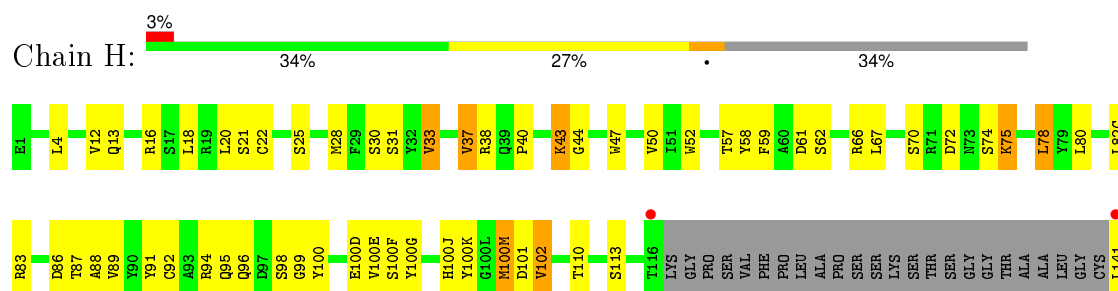
• Molecule 1: Hemagglutinin HA1

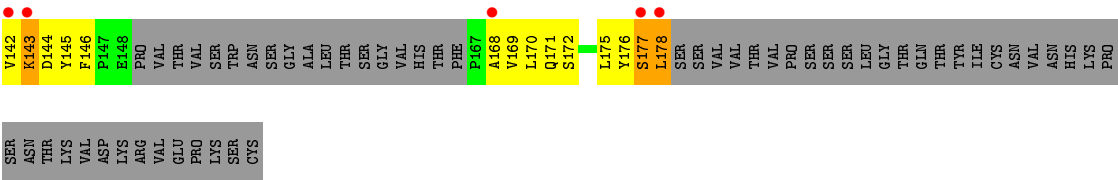


• Molecule 2: Hemagglutinin HA2

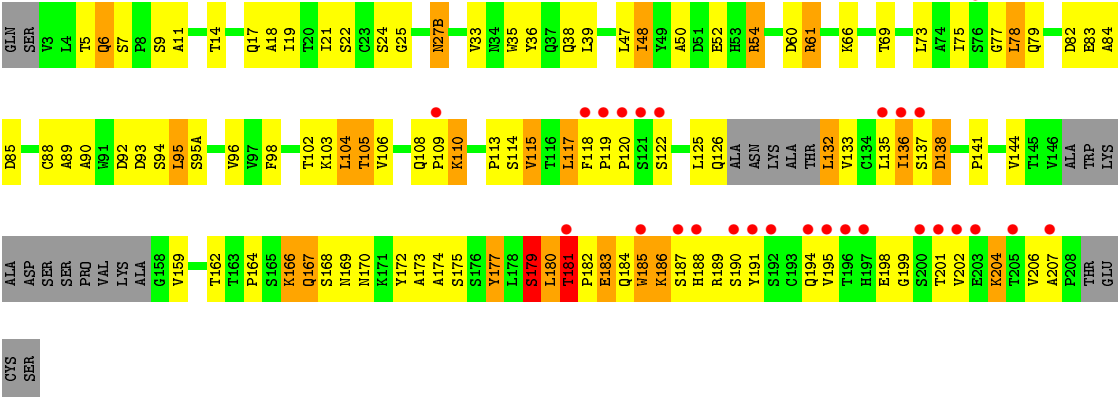


• Molecule 3: Fab 8F8 heavy chain





● Molecule 4: Fab 8F8 light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	136.62Å 136.62Å 142.13Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.72 – 3.00 45.47 – 3.00	Depositor EDS
% Data completeness (in resolution range)	87.8 (44.72-3.00) 94.4 (45.47-3.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.229 , 0.282 0.227 , 0.286	Depositor DCC
R_{free} test set	1499 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	76.8	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 54.7	EDS
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 29407 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6639	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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: 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	0.32	0/2594	0.53	0/3521
2	B	0.34	0/1424	0.60	1/1912 (0.1%)
3	H	0.32	0/1224	0.52	0/1656
4	L	0.32	0/1470	0.69	4/2008 (0.2%)
All	All	0.33	0/6712	0.58	5/9097 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1
4	L	0	5
All	All	0	6

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	147	GLU	N-CA-C	-8.22	88.80	111.00
4	L	181	THR	N-CA-C	-7.27	91.37	111.00
4	L	199	GLY	N-CA-C	-7.10	95.34	113.10
4	L	186	LYS	N-CA-C	-6.38	93.77	111.00
4	L	95	LEU	N-CA-C	-5.76	95.46	111.00

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	146	ASP	Peptide
4	L	179	SER	Peptide
4	L	181	THR	Peptide
4	L	183	GLU	Peptide
4	L	185	TRP	Peptide
4	L	198	GLU	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2534	0	2486	76	0
2	B	1396	0	1299	54	0
3	H	1195	0	1142	44	0
4	L	1436	0	1377	71	0
5	A	50	0	43	1	0
6	A	28	0	25	0	0
All	All	6639	0	6372	231	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 18.

All (231) 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)
2:B:165:GLU:O	2:B:169:ASN:HB2	1.55	1.05
2:B:127:ARG:HD2	2:B:128:ASP:H	1.30	0.96
4:L:187:SER:HA	4:L:190:SER:H	1.33	0.93
4:L:113:PRO:HB2	4:L:136:ILE:HG22	1.56	0.87
3:H:40:PRO:HB2	3:H:43:LYS:HG2	1.57	0.86
1:A:122:PRO:O	1:A:126:ARG:HG2	1.77	0.85
4:L:14:THR:HG22	4:L:17:GLN:HB2	1.61	0.83
2:B:130:VAL:HG12	2:B:140:PHE:HA	1.58	0.82
4:L:187:SER:CB	4:L:188:HIS:HA	2.10	0.82
2:B:143:LYS:HE3	2:B:145:ASP:HA	1.61	0.81
4:L:194:GLN:HA	4:L:202:VAL:O	1.82	0.80
3:H:18:LEU:HD11	3:H:20:LEU:HD13	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:187:SER:HB3	4:L:188:HIS:HA	1.68	0.76
2:B:26:HIS:HD2	2:B:149:MET:HG2	1.52	0.74
3:H:37:VAL:O	3:H:91:TYR:HB2	1.87	0.74
4:L:92:ASP:OD2	4:L:94:SER:HB3	1.88	0.73
4:L:179:SER:O	4:L:180:LEU:HD23	1.88	0.73
3:H:96:GLN:HG3	3:H:100(K):TYR:HB3	1.71	0.73
1:A:137:ARG:HB2	3:H:28:MET:HE1	1.72	0.71
1:A:122:PRO:HD2	1:A:126:ARG:HD2	1.73	0.70
4:L:186:LYS:N	4:L:187:SER:O	2.23	0.69
1:A:211:LYS:HE3	1:A:213:SER:HB2	1.74	0.69
1:A:323:VAL:HB	2:B:12:GLY:HA2	1.73	0.69
1:A:201:TYR:HE2	1:A:248:THR:HG23	1.58	0.68
1:A:164:ALA:O	1:A:246:GLU:HA	1.94	0.67
2:B:127:ARG:HD2	2:B:128:ASP:N	2.08	0.67
4:L:105:THR:HG21	4:L:141:PRO:HB3	1.76	0.67
3:H:13:GLN:O	3:H:16:ARG:HB2	1.96	0.65
1:A:172:SER:HB2	1:A:259:LYS:HD2	1.77	0.65
4:L:138:ASP:HA	4:L:172:TYR:O	1.98	0.64
3:H:94:ARG:HB3	3:H:102:VAL:HG13	1.80	0.63
1:A:120:ILE:HG13	1:A:121:LEU:H	1.64	0.63
1:A:23:THR:O	1:A:25:LYS:HG3	1.99	0.62
1:A:20:ASN:OD1	1:A:22:SER:HB3	2.00	0.61
4:L:185:TRP:O	4:L:186:LYS:HB2	2.00	0.61
2:B:140:PHE:HB3	2:B:142:HIS:O	2.02	0.60
4:L:119:PRO:HB2	4:L:120:PRO:HD2	1.84	0.60
1:A:32:ARG:O	1:A:34:VAL:HG23	2.03	0.59
1:A:278:GLU:HG3	1:A:279:THR:N	2.17	0.59
1:A:208:THR:O	1:A:209:LEU:HB2	2.03	0.59
1:A:209:LEU:HD12	1:A:210:ASN:H	1.68	0.59
2:B:4:GLY:O	2:B:8:GLY:HA3	2.02	0.59
1:A:167:SER:HA	1:A:243:ILE:O	2.03	0.58
2:B:143:LYS:HE3	2:B:145:ASP:CA	2.30	0.58
2:B:147:GLU:C	2:B:147:GLU:OE1	2.42	0.58
2:B:61:THR:O	2:B:62:GLN:HB3	2.04	0.58
3:H:28:MET:HE2	3:H:31:SER:HB3	1.86	0.58
2:B:70:PHE:HB3	2:B:74:GLU:HB2	1.85	0.58
2:B:147:GLU:CD	2:B:147:GLU:C	2.62	0.57
4:L:108:GLN:HB2	4:L:109:PRO:HD2	1.86	0.57
1:A:283:THR:HG22	1:A:301:THR:HG22	1.87	0.57
4:L:187:SER:HA	4:L:190:SER:N	2.13	0.57
2:B:25:HIS:HB3	2:B:34:TYR:CD1	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:59:PHE:CD1	3:H:67:LEU:HD23	2.40	0.57
2:B:107:THR:O	2:B:110:PHE:HB3	2.05	0.57
4:L:114:SER:HB3	4:L:137:SER:HB2	1.85	0.57
4:L:187:SER:CB	4:L:188:HIS:CA	2.83	0.57
1:A:201:TYR:HD2	1:A:248:THR:HG1	1.51	0.57
3:H:67:LEU:HD11	3:H:80:LEU:HD11	1.85	0.57
4:L:179:SER:C	4:L:180:LEU:HD23	2.25	0.57
4:L:6:GLN:NE2	4:L:88:CYS:SG	2.78	0.56
4:L:47:LEU:O	4:L:48:ILE:HD12	2.04	0.56
4:L:167:GLN:HE22	4:L:169:ASN:HB2	1.70	0.56
3:H:143:LYS:HG3	3:H:144:ASP:N	2.20	0.56
1:A:25:LYS:HB3	1:A:33:ASN:ND2	2.20	0.56
3:H:144:ASP:HA	3:H:175:LEU:HG	1.86	0.56
2:B:164:GLU:H	2:B:164:GLU:CD	2.09	0.56
3:H:94:ARG:O	3:H:100(M):MET:HA	2.06	0.56
4:L:18:ALA:HA	4:L:75:ILE:O	2.06	0.56
1:A:15:ILE:HG12	2:B:118:LEU:HD21	1.88	0.55
1:A:71:LEU:O	1:A:148:PHE:HB3	2.06	0.55
1:A:11:ASP:OD2	2:B:143:LYS:HD2	2.06	0.55
4:L:117:LEU:O	4:L:133:VAL:O	2.25	0.55
2:B:151:SER:OG	2:B:157:TYR:HA	2.07	0.55
4:L:204:LYS:O	4:L:206:VAL:HG13	2.07	0.55
4:L:191:TYR:HD2	4:L:207:ALA:HB3	1.72	0.55
4:L:95:LEU:HA	4:L:95(A):SER:C	2.27	0.55
3:H:28:MET:CE	3:H:31:SER:HB3	2.38	0.54
3:H:40:PRO:CB	3:H:43:LYS:HG2	2.33	0.54
1:A:298:HIS:HE1	1:A:300:LEU:HD12	1.71	0.54
3:H:33:VAL:HG22	3:H:95:GLN:HB3	1.89	0.54
2:B:23:GLY:HA3	2:B:36:ALA:HA	1.90	0.54
2:B:26:HIS:CD2	2:B:149:MET:HG2	2.38	0.53
1:A:209:LEU:HD12	1:A:210:ASN:N	2.23	0.53
3:H:38:ARG:NH1	3:H:86:ASP:HA	2.23	0.53
4:L:78:LEU:HD23	4:L:106:VAL:HG12	1.91	0.53
3:H:50:VAL:CG1	3:H:58:TYR:HB2	2.39	0.53
1:A:189:THR:O	1:A:193:THR:HG23	2.09	0.53
4:L:187:SER:HB3	4:L:188:HIS:CA	2.33	0.52
1:A:45:LYS:HG2	1:A:296:ASN:HD21	1.75	0.52
1:A:134:GLY:HA3	3:H:100:TYR:OH	2.08	0.52
1:A:120:ILE:HG13	1:A:121:LEU:N	2.25	0.52
3:H:144:ASP:H	3:H:177:SER:HB3	1.74	0.52
3:H:50:VAL:HG12	3:H:58:TYR:HB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:47:TRP:CG	4:L:96:VAL:HB	2.45	0.52
4:L:11:ALA:O	4:L:104:LEU:HA	2.10	0.52
1:A:50:LYS:HB3	1:A:275:GLU:HG3	1.93	0.51
4:L:6:GLN:HE22	4:L:88:CYS:H	1.56	0.51
1:A:206:THR:HG22	1:A:243:ILE:HA	1.92	0.51
4:L:108:GLN:HB2	4:L:109:PRO:CD	2.40	0.51
1:A:17:TYR:CE2	2:B:6:ILE:HG12	2.46	0.51
3:H:100(J):HIS:O	4:L:50:ALA:HB2	2.11	0.51
3:H:141:LEU:HG	3:H:142:VAL:N	2.26	0.51
1:A:42:ILE:HG13	1:A:314:LEU:HB3	1.93	0.51
4:L:185:TRP:N	4:L:185:TRP:CD1	2.79	0.50
1:A:226:GLN:OE1	3:H:99:GLY:HA3	2.11	0.50
3:H:96:GLN:HB3	3:H:101:ASP:OD2	2.12	0.50
1:A:307:LYS:HG3	2:B:62:GLN:HG2	1.93	0.50
1:A:224:ASN:O	1:A:226:GLN:HG2	2.11	0.50
4:L:115:VAL:HB	4:L:136:ILE:HG23	1.94	0.49
2:B:28:ASN:HD21	2:B:145:ASP:C	2.15	0.49
4:L:132:LEU:HD11	4:L:179:SER:O	2.12	0.49
1:A:220:ARG:HD2	1:A:227:GLY:O	2.12	0.49
3:H:43:LYS:HG3	3:H:44:GLY:N	2.26	0.49
4:L:166:LYS:NZ	4:L:170:ASN:HA	2.27	0.49
2:B:55:VAL:HG23	2:B:56:ILE:N	2.28	0.49
4:L:166:LYS:NZ	4:L:167:GLN:H	2.11	0.48
4:L:181:THR:O	4:L:183:GLU:N	2.43	0.48
1:A:137:ARG:CB	3:H:28:MET:HE1	2.43	0.48
1:A:25:LYS:HD2	1:A:33:ASN:OD1	2.13	0.48
4:L:114:SER:HB3	4:L:137:SER:CB	2.42	0.48
3:H:143:LYS:HA	3:H:177:SER:HB3	1.95	0.48
4:L:95:LEU:HG	4:L:95(A):SER:HA	1.95	0.48
1:A:45:LYS:HA	1:A:296:ASN:OD1	2.13	0.48
2:B:71:SER:OG	2:B:74:GLU:HG3	2.13	0.48
2:B:129:ASN:O	2:B:141:TYR:HB2	2.14	0.48
1:A:43:LEU:HB2	1:A:314:LEU:HB2	1.96	0.48
1:A:316:LEU:HD11	2:B:55:VAL:HG21	1.94	0.48
4:L:82:ASP:O	4:L:104:LEU:HD21	2.14	0.48
1:A:188:GLU:H	1:A:188:GLU:CD	2.16	0.48
1:A:89:GLU:OE1	1:A:269:LYS:HE2	2.14	0.48
4:L:180:LEU:HB3	4:L:182:PRO:HA	1.95	0.48
2:B:151:SER:HA	2:B:154:ASN:ND2	2.29	0.47
1:A:186:ASN:OD1	1:A:227:GLY:HA3	2.14	0.47
1:A:295:HIS:HD2	1:A:306:PRO:HG2	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:PRO:O	1:A:294:PHE:CG	2.67	0.47
1:A:299:PRO:HG3	1:A:308:TYR:CE2	2.50	0.47
2:B:71:SER:H	2:B:74:GLU:HG3	1.80	0.47
3:H:66:ARG:HH22	3:H:86:ASP:CG	2.18	0.47
3:H:21:SER:HA	3:H:78:LEU:O	2.15	0.47
1:A:325:GLN:HB2	2:B:15:GLN:OE1	2.15	0.47
2:B:55:VAL:HG23	2:B:56:ILE:H	1.81	0.46
4:L:185:TRP:O	4:L:186:LYS:CB	2.63	0.46
4:L:19:ILE:HA	4:L:19:ILE:HD12	1.76	0.46
1:A:67:ALA:O	1:A:71:LEU:HB2	2.15	0.46
2:B:56:ILE:HG13	2:B:57:GLU:N	2.30	0.46
4:L:186:LYS:HA	4:L:190:SER:N	2.30	0.46
1:A:25:LYS:CB	1:A:33:ASN:HD21	2.28	0.46
2:B:52:VAL:HA	2:B:55:VAL:HG22	1.97	0.46
2:B:127:ARG:CD	2:B:128:ASP:H	2.16	0.46
4:L:25:GLY:O	4:L:69:THR:HB	2.17	0.45
1:A:71:LEU:HA	1:A:71:LEU:HD12	1.83	0.45
3:H:168:ALA:HA	3:H:178:LEU:HB3	1.99	0.45
4:L:113:PRO:HA	4:L:138:ASP:O	2.16	0.45
3:H:43:LYS:CG	3:H:44:GLY:N	2.80	0.45
2:B:125:GLN:HE22	2:B:157:TYR:H	1.63	0.45
4:L:136:ILE:HG12	4:L:174:ALA:HB3	1.99	0.45
1:A:18:HIS:ND1	1:A:19:ALA:N	2.64	0.45
1:A:180:TRP:CZ2	1:A:233:SER:HB2	2.51	0.44
2:B:125:GLN:HB3	2:B:126:LEU:HD23	1.99	0.44
1:A:320:LEU:HB3	2:B:111:HIS:CD2	2.52	0.44
4:L:180:LEU:HD22	4:L:180:LEU:HA	1.81	0.44
4:L:61:ARG:HD3	4:L:77:GLY:O	2.17	0.44
3:H:87:THR:O	3:H:88:ALA:HB2	2.18	0.44
1:A:54:ASN:OD1	1:A:280:LYS:HE3	2.17	0.44
1:A:119:LYS:HE2	1:A:255:GLU:OE2	2.18	0.44
1:A:52:CYS:HB3	1:A:277:CYS:O	2.18	0.44
2:B:140:PHE:HE2	2:B:149:MET:HE1	1.83	0.44
3:H:4:LEU:HB3	3:H:22:CYS:SG	2.58	0.44
4:L:36:TYR:HE1	4:L:89:ALA:HB3	1.82	0.44
2:B:148:CYS:HA	2:B:151:SER:HB3	2.00	0.43
1:A:25:LYS:HB3	1:A:33:ASN:HD21	1.81	0.43
3:H:83:ARG:HG2	3:H:86:ASP:OD2	2.18	0.43
4:L:186:LYS:HE2	4:L:189:ARG:HG3	2.00	0.43
1:A:204:VAL:HA	1:A:244:ASN:O	2.18	0.43
1:A:32:ARG:HD3	1:A:32:ARG:HA	1.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:83:GLU:OE1	4:L:106:VAL:HG22	2.18	0.43
1:A:214:THR:HA	1:A:215:PRO:HD3	1.71	0.43
4:L:119:PRO:CB	4:L:120:PRO:HD2	2.48	0.43
3:H:72:ASP:OD1	3:H:75:LYS:HB2	2.17	0.43
2:B:37:ASP:O	2:B:41:THR:HB	2.18	0.43
1:A:159:SER:O	1:A:196:GLN:HG3	2.18	0.43
2:B:164:GLU:HA	2:B:167:LYS:HD3	2.00	0.43
1:A:58:PRO:HB3	1:A:86:TYR:CZ	2.54	0.43
4:L:92:ASP:O	4:L:93:ASP:HB2	2.19	0.43
1:A:316:LEU:HD23	1:A:316:LEU:HA	1.89	0.42
2:B:14:TRP:HB2	2:B:17:MET:HG2	2.00	0.42
3:H:78:LEU:HA	3:H:78:LEU:HD23	1.71	0.42
1:A:129:GLN:HB3	1:A:162:PRO:HG3	2.02	0.42
1:A:116(C):GLU:HG2	1:A:261:SER:OG	2.19	0.42
4:L:186:LYS:CE	4:L:189:ARG:HG3	2.50	0.42
2:B:126:LEU:HD23	2:B:126:LEU:N	2.35	0.42
1:A:222:LYS:N	1:A:222:LYS:HD2	2.34	0.42
4:L:38:GLN:O	4:L:84:ALA:HB1	2.20	0.42
4:L:164:PRO:HA	4:L:173:ALA:O	2.19	0.42
2:B:149:MET:O	2:B:152:VAL:HB	2.19	0.42
4:L:166:LYS:HZ2	4:L:166:LYS:HA	1.85	0.42
4:L:117:LEU:HD22	4:L:117:LEU:HA	1.76	0.42
4:L:177:TYR:CD1	4:L:177:TYR:N	2.87	0.42
4:L:186:LYS:HG2	4:L:189:ARG:HB3	2.02	0.41
1:A:120:ILE:O	1:A:122:PRO:HD3	2.20	0.41
1:A:203:SER:O	1:A:245:PHE:HA	2.21	0.41
3:H:40:PRO:HD2	3:H:44:GLY:O	2.20	0.41
4:L:33:VAL:HG22	4:L:90:ALA:HB2	2.02	0.41
5:A:403:BMA:H62	5:A:404:MAN:H2	1.66	0.41
4:L:35:TRP:CE3	4:L:73:LEU:HD22	2.55	0.41
1:A:314:LEU:HD23	1:A:314:LEU:HA	1.81	0.41
1:A:223:VAL:O	1:A:224:ASN:HB2	2.20	0.41
3:H:22:CYS:HB3	3:H:78:LEU:HB2	2.02	0.41
3:H:100(D):GLU:HB2	3:H:100(G):TYR:CE2	2.56	0.41
2:B:126:LEU:HD21	2:B:152:VAL:HG13	2.02	0.41
1:A:295:HIS:CD2	1:A:306:PRO:HG2	2.55	0.41
4:L:183:GLU:O	4:L:185:TRP:N	2.45	0.41
2:B:71:SER:CB	2:B:74:GLU:HG3	2.51	0.41
2:B:167:LYS:HG2	2:B:168:LEU:N	2.35	0.41
4:L:78:LEU:CD2	4:L:106:VAL:HG12	2.50	0.41
3:H:47:TRP:HB2	4:L:98:PHE:HE1	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:LYS:O	1:A:286:GLY:HA2	2.21	0.41
1:A:202:VAL:HA	1:A:247:SER:HB2	2.01	0.41
2:B:54:SER:O	2:B:58:LYS:HB2	2.20	0.41
2:B:3:PHE:HD2	2:B:112:ASP:HB3	1.86	0.41
4:L:54:ARG:NH1	4:L:60:ASP:HA	2.36	0.41
2:B:142:HIS:CD2	2:B:162:TYR:HB3	2.56	0.41
1:A:42:ILE:CG1	1:A:314:LEU:HB3	2.51	0.41
4:L:110:LYS:HG2	4:L:110:LYS:H	1.53	0.41
1:A:58:PRO:HB3	1:A:86:TYR:CE2	2.56	0.40
3:H:168:ALA:HA	3:H:178:LEU:CB	2.52	0.40
3:H:145:TYR:HD1	3:H:176:TYR:O	2.05	0.40
2:B:165:GLU:O	2:B:169:ASN:CB	2.46	0.40
2:B:57:GLU:HA	2:B:60:ASN:OD1	2.22	0.40
4:L:85:ASP:HA	4:L:102:THR:O	2.22	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	322/327 (98%)	303 (94%)	19 (6%)	0	100	100
2	B	170/174 (98%)	154 (91%)	15 (9%)	1 (1%)	30	72
3	H	147/233 (63%)	136 (92%)	11 (8%)	0	100	100
4	L	190/218 (87%)	157 (83%)	30 (16%)	3 (2%)	12	48
All	All	829/952 (87%)	750 (90%)	75 (9%)	4 (0%)	34	76

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	L	184	GLN

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Mol	Chain	Res	Type
4	L	27(B)	ASN
2	B	68	LYS
4	L	179	SER

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	285/288 (99%)	246 (86%)	39 (14%)	4	20
2	B	149/151 (99%)	118 (79%)	31 (21%)	1	7
3	H	128/197 (65%)	96 (75%)	32 (25%)	1	3
4	L	162/179 (90%)	119 (74%)	43 (26%)	0	3
All	All	724/815 (89%)	579 (80%)	145 (20%)	1	8

All (145) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ILE
1	A	27	ASP
1	A	30	LEU
1	A	32	ARG
1	A	36	VAL
1	A	45	LYS
1	A	51	LEU
1	A	63	ASP
1	A	71	LEU
1	A	78	ARG
1	A	80	LEU
1	A	97	CYS
1	A	101	SER
1	A	133	THR
1	A	156	GLU
1	A	167	SER
1	A	178	ILE
1	A	182	VAL

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Mol	Chain	Res	Type
1	A	186	ASN
1	A	193	THR
1	A	201	TYR
1	A	203	SER
1	A	208	THR
1	A	214	THR
1	A	222	LYS
1	A	235	THR
1	A	243	ILE
1	A	247	SER
1	A	261	SER
1	A	264	SER
1	A	270	THR
1	A	274	LEU
1	A	275	GLU
1	A	278	GLU
1	A	291	THR
1	A	315	VAL
1	A	320	LEU
1	A	323	VAL
1	A	325	GLN
2	B	10	ILE
2	B	17	MET
2	B	22	TYR
2	B	24	TYR
2	B	25	HIS
2	B	26	HIS
2	B	41	THR
2	B	49	THR
2	B	56	ILE
2	B	58	LYS
2	B	59	MET
2	B	61	THR
2	B	62	GLN
2	B	73	LEU
2	B	75	ARG
2	B	103	GLU
2	B	113	SER
2	B	116	LYS
2	B	118	LEU
2	B	121	LYS
2	B	123	ARG

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Mol	Chain	Res	Type
2	B	125	GLN
2	B	126	LEU
2	B	127	ARG
2	B	145	ASP
2	B	147	GLU
2	B	149	MET
2	B	156	THR
2	B	157	TYR
2	B	166	SER
2	B	167	LYS
3	H	12	VAL
3	H	25	SER
3	H	30	SER
3	H	33	VAL
3	H	37	VAL
3	H	43	LYS
3	H	52	TRP
3	H	57	THR
3	H	61	ASP
3	H	62	SER
3	H	70	SER
3	H	74	SER
3	H	75	LYS
3	H	78	LEU
3	H	82(C)	LEU
3	H	89	VAL
3	H	92	CYS
3	H	98	SER
3	H	100(E)	VAL
3	H	100(F)	SER
3	H	100(M)	MET
3	H	102	VAL
3	H	110	THR
3	H	113	SER
3	H	143	LYS
3	H	146	PHE
3	H	169	VAL
3	H	170	LEU
3	H	171	GLN
3	H	172	SER
3	H	177	SER
3	H	178	LEU

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Mol	Chain	Res	Type
4	L	5	THR
4	L	6	GLN
4	L	7	SER
4	L	9	SER
4	L	21	ILE
4	L	22	SER
4	L	24	SER
4	L	27(B)	ASN
4	L	39	LEU
4	L	48	ILE
4	L	52	GLU
4	L	54	ARG
4	L	61	ARG
4	L	66	LYS
4	L	78	LEU
4	L	79	GLN
4	L	103	LYS
4	L	104	LEU
4	L	105	THR
4	L	110	LYS
4	L	115	VAL
4	L	117	LEU
4	L	118	PHE
4	L	122	SER
4	L	125	LEU
4	L	126	GLN
4	L	132	LEU
4	L	135	LEU
4	L	136	ILE
4	L	138	ASP
4	L	144	VAL
4	L	159	VAL
4	L	162	THR
4	L	166	LYS
4	L	167	GLN
4	L	168	SER
4	L	175	SER
4	L	177	TYR
4	L	180	LEU
4	L	181	THR
4	L	195	VAL
4	L	201	THR

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Mol	Chain	Res	Type
4	L	204	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	HIS
1	A	295	HIS
2	B	125	GLN
2	B	129	ASN
2	B	142	HIS
3	H	96	GLN
4	L	167	GLN
4	L	170	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 ⓘ

6 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	401	1,5	14,14,15	0.45	0	15,19,21	1.02	1 (6%)
5	NAG	A	402	5	14,14,15	0.56	0	15,19,21	0.73	0
5	BMA	A	403	5	11,11,12	0.57	0	14,15,17	0.60	0
5	MAN	A	404	5	11,11,12	0.58	0	14,15,17	0.75	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	405	1,6	14,14,15	0.47	0	15,19,21	0.96	0
6	NAG	A	406	6	14,14,15	0.48	0	15,19,21	0.64	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	401	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	402	5	-	0/6/23/26	0/1/1/1
5	BMA	A	403	5	-	0/2/19/22	0/1/1/1
5	MAN	A	404	5	-	0/2/19/22	0/1/1/1
6	NAG	A	405	1,6	-	0/6/23/26	0/1/1/1
6	NAG	A	406	6	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	401	NAG	C3-C4-C5	2.18	114.00	110.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	403	BMA	1	0
5	A	404	MAN	1	0

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	324/327 (99%)	-0.29	1 (0%) 94 84	49, 71, 118, 148	0
2	B	172/174 (98%)	0.10	2 (1%) 81 55	48, 113, 161, 209	0
3	H	153/233 (65%)	-0.07	7 (4%) 36 14	49, 80, 161, 176	0
4	L	196/218 (89%)	0.52	27 (13%) 4 1	63, 130, 187, 218	0
All	All	845/952 (88%)	0.02	37 (4%) 38 16	48, 87, 165, 218	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	L	188	HIS	5.4
3	H	142	VAL	5.0
4	L	136	ILE	4.4
4	L	200	SER	4.0
4	L	194	GLN	4.0
4	L	120	PRO	3.7
4	L	135	LEU	3.6
4	L	191	TYR	3.6
4	L	201	THR	3.6
4	L	119	PRO	3.5
4	L	203	GLU	3.3
3	H	178	LEU	3.3
3	H	177	SER	3.2
4	L	137	SER	3.1
1	A	12	GLN	2.8
4	L	197	HIS	2.7
2	B	16	GLY	2.7
3	H	143	LYS	2.6
4	L	196	THR	2.6
3	H	116	THR	2.5
4	L	195	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
4	L	122	SER	2.5
4	L	118	PHE	2.4
4	L	187	SER	2.4
4	L	185	TRP	2.3
4	L	205	THR	2.2
4	L	202	VAL	2.2
4	L	207	ALA	2.2
4	L	192	SER	2.2
3	H	141	LEU	2.2
4	L	121	SER	2.2
4	L	76	SER	2.1
2	B	19	ASP	2.1
3	H	168	ALA	2.1
4	L	190	SER	2.0
4	L	109	PRO	2.0
4	L	181	THR	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(\AA^2)	Q<0.9
5	MAN	A	404	11/12	0.92	0.33	-	140,147,148,148	0
6	NAG	A	406	14/15	0.83	0.33	-	125,135,137,138	0
5	NAG	A	401	14/15	0.94	0.14	-	80,92,99,106	0
5	BMA	A	403	11/12	0.83	0.23	-	140,144,147,147	0
5	NAG	A	402	14/15	0.85	0.16	-	110,119,126,134	0
6	NAG	A	405	14/15	0.89	0.22	-	100,108,118,126	0

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.