



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 3, 2020 – 04:04 PM EST

PDB ID : 6OCB
Title : Crystal structure of FluA-20 Fab in complex with the head domain of H3
(A/Hong Kong/1/1968)
Authors : Wilson, I.A.; Lang, S.
Deposited on : 2019-03-22
Resolution : 2.10 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.6
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.14.6

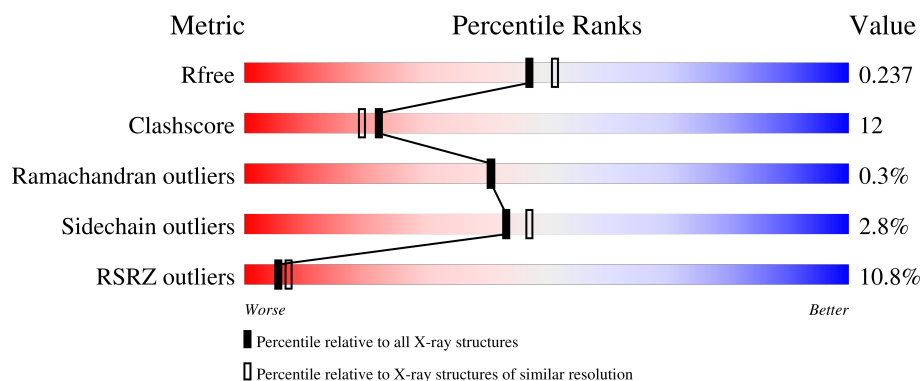
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 respectively. 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	H	235	<div> <div>11%</div> <div> <div></div> <div>73%</div> <div>19%</div> <div>6%</div> </div> </div>
2	L	214	<div> <div>6%</div> <div> <div></div> <div>83%</div> <div>14%</div> <div>..</div> </div> </div>
3	A	274	<div> <div>14%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>..</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5746 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 Heavy chain of FluA-20 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	222	Total	C	N	O	S	0	0	0
			1662	1050	273	332	7			

- Molecule 2 is a protein called Light chain of FluA-20 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	213	Total	C	N	O	S	0	0	0
			1646	1031	281	329	5			

- Molecule 3 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	267	Total	C	N	O	S	0	1	0
			2079	1306	364	398	11			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	310	GLY	-	expression tag	UNP Q91MA7
A	311	HIS	-	expression tag	UNP Q91MA7
A	312	HIS	-	expression tag	UNP Q91MA7
A	313	HIS	-	expression tag	UNP Q91MA7
A	314	HIS	-	expression tag	UNP Q91MA7
A	315	HIS	-	expression tag	UNP Q91MA7
A	316	HIS	-	expression tag	UNP Q91MA7

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

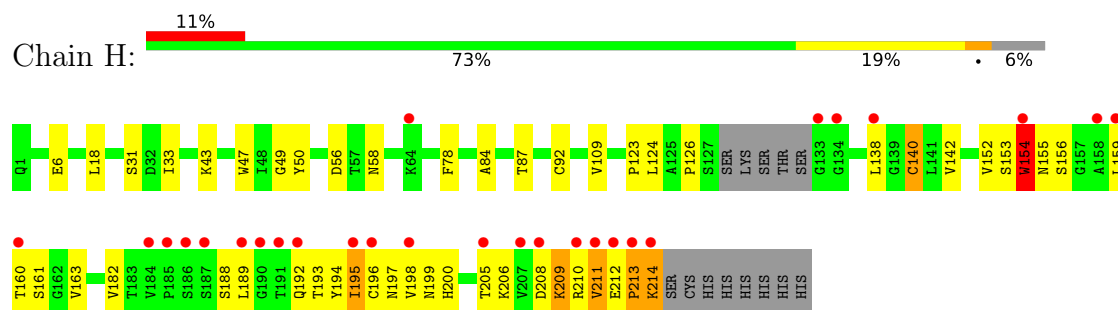
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	76	Total	O	0	0
			76	76		
5	L	114	Total	O	0	0
			114	114		
5	A	155	Total	O	0	0
			155	155		

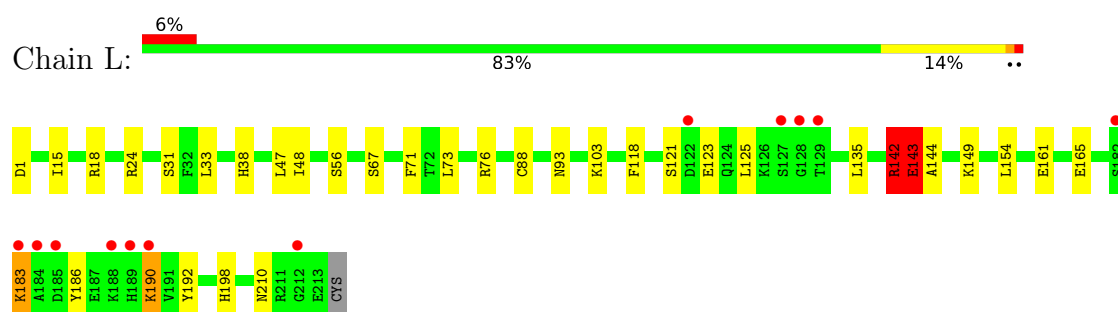
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

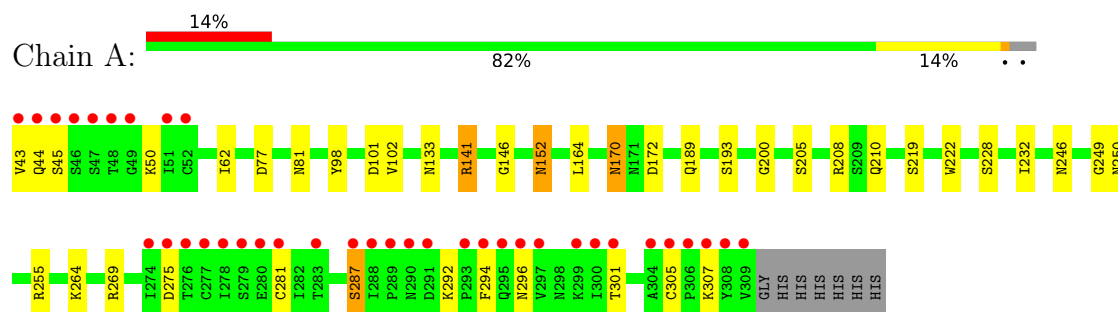
• Molecule 1: Heavy chain of FluA-20 Fab



• Molecule 2: Light chain of FluA-20 Fab



• Molecule 3: Hemagglutinin



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	84.55Å 84.55Å 271.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.81 – 2.10 44.88 – 2.09	Depositor EDS
% Data completeness (in resolution range)	99.3 (37.81-2.10) 99.4 (44.88-2.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.08Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.202 , 0.239 0.203 , 0.237	Depositor DCC
R_{free} test set	2965 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	33.1	Xtriage
Anisotropy	0.511	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5746	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section:
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	H	0.70	1/1704 (0.1%)	0.91	3/2330 (0.1%)
2	L	0.76	3/1683 (0.2%)	0.96	7/2285 (0.3%)
3	A	0.69	0/2131	0.81	2/2903 (0.1%)
All	All	0.71	4/5518 (0.1%)	0.89	12/7518 (0.2%)

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
1	H	0	1
2	L	0	1
3	A	0	1
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	143	GLU	CB-CG	-12.88	1.27	1.52
2	L	143	GLU	CG-CD	-8.96	1.38	1.51
1	H	140	CYS	CB-SG	-6.65	1.71	1.82
2	L	93	ASN	C-N	-5.49	1.21	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	24	ARG	NE-CZ-NH1	9.11	124.86	120.30
2	L	144	ALA	N-CA-C	-7.71	90.18	111.00
1	H	140	CYS	CA-CB-SG	6.99	126.58	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	143	GLU	CB-CA-C	-6.72	96.95	110.40
3	A	141	ARG	CG-CD-NE	-6.16	98.86	111.80
2	L	142	ARG	CG-CD-NE	6.04	124.49	111.80
2	L	24	ARG	NE-CZ-NH2	-6.01	117.29	120.30
2	L	47	LEU	CA-CB-CG	5.80	128.63	115.30
1	H	138	LEU	CB-CG-CD2	5.71	120.72	111.00
1	H	154	TRP	O-C-N	-5.62	113.71	122.70
2	L	190	LYS	CG-CD-CE	-5.32	95.95	111.90
3	A	208	ARG	N-CA-CB	-5.14	101.35	110.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	219	SER	Mainchain
1	H	154	TRP	Mainchain
2	L	143	GLU	Mainchain

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	H	1662	0	1618	49	0
2	L	1646	0	1608	47	0
3	A	2079	0	2018	31	0
4	A	14	0	13	0	0
5	A	155	0	0	3	0
5	H	76	0	0	5	0
5	L	114	0	0	5	0
All	All	5746	0	5257	125	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 (125) 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:L:125:LEU:HB3	2:L:183:LYS:NZ	1.14	1.43
2:L:125:LEU:CB	2:L:183:LYS:HZ1	1.60	1.15
1:H:212:GLU:HG3	1:H:213:PRO:HD2	1.29	1.14
2:L:125:LEU:CB	2:L:183:LYS:NZ	2.09	1.13
3:A:43:VAL:N	3:A:292:LYS:HZ2	1.58	1.01
2:L:142:ARG:NH1	2:L:143:GLU:OE2	1.95	0.98
1:H:212:GLU:CG	1:H:213:PRO:HD2	1.93	0.97
1:H:214:LYS:HD3	1:H:214:LYS:H	1.28	0.96
2:L:125:LEU:HB3	2:L:183:LYS:HZ3	1.17	0.93
2:L:161:GLU:OE1	5:L:301:HOH:O	1.87	0.91
1:H:154:TRP:O	1:H:155:ASN:HB2	1.70	0.90
1:H:159:LEU:HD22	1:H:194:TYR:CE1	2.09	0.88
1:H:56:ASP:OD2	5:H:302:HOH:O	1.92	0.85
1:H:188:SER:HB3	5:H:303:HOH:O	1.77	0.85
3:A:101:ASP:OD1	5:A:801:HOH:O	1.96	0.83
1:H:212:GLU:OE1	1:H:213:PRO:HD3	1.78	0.82
2:L:190:LYS:O	2:L:190:LYS:NZ	2.17	0.77
1:H:159:LEU:HD22	1:H:194:TYR:HE1	1.50	0.76
1:H:153:SER:HB2	1:H:197:ASN:ND2	2.01	0.76
1:H:188:SER:O	5:H:303:HOH:O	2.03	0.75
2:L:143:GLU:OE2	2:L:143:GLU:N	2.21	0.73
2:L:190:LYS:HZ1	2:L:210:ASN:HA	1.51	0.73
2:L:125:LEU:O	2:L:183:LYS:HE2	1.89	0.72
2:L:190:LYS:C	2:L:190:LYS:HZ2	1.92	0.72
2:L:142:ARG:HD3	2:L:143:GLU:OE2	1.90	0.71
2:L:190:LYS:HE2	2:L:210:ASN:HB3	1.73	0.70
1:H:208:ASP:O	1:H:209:LYS:HD2	1.91	0.70
1:H:214:LYS:CD	1:H:214:LYS:H	2.01	0.69
2:L:125:LEU:CB	2:L:183:LYS:HZ3	1.89	0.69
2:L:142:ARG:NE	5:L:302:HOH:O	2.05	0.68
2:L:123:GLU:N	2:L:123:GLU:OE1	2.25	0.68
3:A:133:ASN:H	3:A:152:ASN:HD21	1.39	0.68
1:H:212:GLU:OE1	1:H:213:PRO:CD	2.41	0.67
1:H:209:LYS:N	5:H:301:HOH:O	2.27	0.67
2:L:190:LYS:C	2:L:190:LYS:NZ	2.48	0.66
1:H:195:ILE:HG12	1:H:195:ILE:O	1.94	0.66
3:A:133:ASN:H	3:A:152:ASN:ND2	1.94	0.65
3:A:141:ARG:HG3	3:A:146:GLY:HA3	1.77	0.65
3:A:43:VAL:N	3:A:292:LYS:NZ	2.41	0.64
1:H:205:THR:O	1:H:206:LYS:HD3	1.98	0.63
2:L:18:ARG:HH21	2:L:76:ARG:NH2	1.97	0.63
2:L:190:LYS:NZ	2:L:210:ASN:HA	2.13	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:159:LEU:HD22	1:H:194:TYR:CD1	2.33	0.62
1:H:84:ALA:O	1:H:87:THR:HG22	1.98	0.62
1:H:212:GLU:CG	1:H:213:PRO:CD	2.76	0.61
1:H:210:ARG:HH11	1:H:210:ARG:HG3	1.67	0.60
1:H:18:LEU:HD13	1:H:109:VAL:HG11	1.85	0.59
1:H:6:GLU:HG3	1:H:92:CYS:SG	2.42	0.58
3:A:281:CYS:O	3:A:287:SER:OG	2.19	0.58
3:A:189:GLN:NE2	3:A:193:SER:OG	2.37	0.58
1:H:153:SER:HB2	1:H:197:ASN:HD21	1.67	0.57
2:L:143:GLU:O	2:L:198:HIS:HD2	1.87	0.57
2:L:125:LEU:HB3	2:L:183:LYS:HZ1	0.73	0.56
1:H:152:VAL:HG22	1:H:198:VAL:HG22	1.87	0.55
3:A:301:THR:OG1	3:A:305:CYS:SG	2.61	0.55
1:H:142:VAL:HG22	1:H:198:VAL:HG21	1.88	0.54
1:H:154:TRP:O	1:H:155:ASN:CB	2.35	0.54
2:L:56:SER:HB2	3:A:222:TRP:CZ2	2.42	0.54
3:A:269:ARG:NH1	5:A:805:HOH:O	2.40	0.53
1:H:163:VAL:HG22	1:H:182:VAL:CG1	2.38	0.53
3:A:98:TYR:HH	3:A:228[B]:SER:HB2	1.74	0.53
2:L:103:LYS:NZ	2:L:165:GLU:OE2	2.38	0.53
1:H:47:TRP:CH2	1:H:49:GLY:HA2	2.44	0.53
2:L:121:SER:O	2:L:125:LEU:HD13	2.08	0.52
3:A:81:ASN:ND2	5:A:807:HOH:O	2.41	0.52
1:H:163:VAL:HG22	1:H:182:VAL:HG12	1.92	0.52
1:H:210:ARG:NH1	1:H:210:ARG:HG3	2.24	0.52
2:L:142:ARG:CZ	2:L:143:GLU:OE1	2.58	0.52
3:A:50:LYS:HD3	3:A:275:ASP:OD2	2.09	0.51
1:H:154:TRP:CH2	1:H:196:CYS:HB3	2.46	0.51
3:A:77:ASP:OD2	3:A:141:ARG:NH2	2.31	0.51
3:A:133:ASN:ND2	3:A:255:ARG:HH22	2.08	0.51
2:L:38:HIS:HD2	5:L:303:HOH:O	1.93	0.50
2:L:142:ARG:HD3	2:L:143:GLU:H	1.76	0.50
3:A:45:SER:HA	3:A:296:ASN:OD1	2.12	0.50
1:H:159:LEU:CD2	1:H:194:TYR:CE1	2.90	0.50
2:L:18:ARG:HE	2:L:76:ARG:NH1	2.09	0.49
2:L:186:TYR:O	2:L:192:TYR:OH	2.30	0.49
1:H:208:ASP:C	1:H:209:LYS:HD2	2.32	0.49
2:L:142:ARG:CZ	2:L:143:GLU:OE2	2.57	0.49
2:L:143:GLU:O	2:L:198:HIS:CD2	2.65	0.49
2:L:142:ARG:NH1	2:L:143:GLU:CD	2.65	0.49
2:L:103:LYS:CE	2:L:165:GLU:OE2	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:102:VAL:HG22	3:A:232:ILE:HB	1.95	0.48
1:H:200:HIS:HB3	1:H:205:THR:OG1	2.13	0.48
1:H:126:PRO:HG3	1:H:189:LEU:HD22	1.96	0.48
1:H:50:TYR:CE2	1:H:58:ASN:HB3	2.49	0.48
3:A:264:LYS:HG2	3:A:264:LYS:O	2.14	0.48
1:H:196:CYS:N	5:H:301:HOH:O	2.46	0.47
1:H:199:ASN:CG	1:H:206:LYS:HD2	2.34	0.47
3:A:205:SER:HB3	3:A:210:GLN:HG3	1.96	0.47
3:A:44:GLN:N	3:A:292:LYS:HD3	2.30	0.46
3:A:164:LEU:O	3:A:246:ASN:HA	2.15	0.46
1:H:154:TRP:O	1:H:155:ASN:C	2.52	0.46
2:L:67:SER:HA	2:L:71:PHE:CE2	2.51	0.46
3:A:170:ASN:ND2	3:A:172:ASP:H	2.14	0.46
2:L:142:ARG:HD2	2:L:142:ARG:H	1.81	0.46
1:H:124:LEU:HB3	2:L:118:PHE:CD1	2.50	0.45
3:A:133:ASN:HD22	3:A:255:ARG:HH22	1.64	0.45
1:H:212:GLU:CD	1:H:213:PRO:CD	2.86	0.44
2:L:31:SER:OG	5:L:304:HOH:O	2.21	0.44
2:L:33:LEU:HD21	2:L:88:CYS:HB2	2.00	0.44
2:L:48:ILE:HD12	2:L:73:LEU:HD13	1.99	0.43
2:L:142:ARG:CZ	2:L:143:GLU:CD	2.87	0.43
3:A:294:PHE:HA	3:A:307:LYS:O	2.18	0.43
1:H:123:PRO:HB3	1:H:211:VAL:HG13	2.00	0.42
1:H:78:PHE:CZ	1:H:92:CYS:HB2	2.54	0.42
3:A:170:ASN:HD22	3:A:170:ASN:C	2.21	0.42
3:A:200:GLY:HA3	3:A:250:ASN:ND2	2.35	0.42
2:L:142:ARG:HD3	2:L:143:GLU:N	2.34	0.42
2:L:142:ARG:CD	2:L:142:ARG:H	2.32	0.42
3:A:170:ASN:HD21	3:A:172:ASP:HB2	1.85	0.42
1:H:212:GLU:CD	1:H:213:PRO:HD2	2.40	0.41
3:A:170:ASN:HD22	3:A:172:ASP:H	1.67	0.41
3:A:249:GLY:C	3:A:250:ASN:HD22	2.24	0.41
2:L:15:ILE:HD12	5:L:317:HOH:O	2.20	0.41
1:H:43:LYS:HE2	1:H:43:LYS:HB3	1.57	0.41
2:L:149:LYS:HG2	2:L:154:LEU:HD13	2.03	0.41
2:L:18:ARG:HH21	2:L:76:ARG:CZ	2.34	0.41
1:H:142:VAL:CG2	1:H:198:VAL:HG21	2.51	0.41
1:H:160:THR:HG23	1:H:161:SER:N	2.36	0.40
3:A:50:LYS:CD	3:A:275:ASP:OD2	2.69	0.40
1:H:214:LYS:N	1:H:214:LYS:CD	2.75	0.40
2:L:123:GLU:CD	2:L:123:GLU:H	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:142:ARG:N	2:L:142:ARG:CD	2.84	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	H	218/235 (93%)	209 (96%)	8 (4%)	1 (0%)	29	26
2	L	211/214 (99%)	207 (98%)	4 (2%)	0	100	100
3	A	266/274 (97%)	258 (97%)	7 (3%)	1 (0%)	34	32
All	All	695/723 (96%)	674 (97%)	19 (3%)	2 (0%)	41	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	62	ILE
1	H	213	PRO

5.3.2 Protein sidechains [i](#)

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	H	192/205 (94%)	182 (95%)	10 (5%)	23	21
2	L	189/190 (100%)	185 (98%)	4 (2%)	53	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	A	237/242 (98%)	234 (99%)	3 (1%)	69	75
All	All	618/637 (97%)	601 (97%)	17 (3%)	43	47

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	31	SER
1	H	33	ILE
1	H	140	CYS
1	H	156	SER
1	H	192	GLN
1	H	193	THR
1	H	195	ILE
1	H	209	LYS
1	H	211	VAL
1	H	214	LYS
2	L	1	ASP
2	L	135	LEU
2	L	142	ARG
2	L	183	LYS
3	A	152	ASN
3	A	170	ASN
3	A	287	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	L	38	HIS
2	L	199	GLN
3	A	132	GLN
3	A	133	ASN
3	A	152	ASN
3	A	170	ASN
3	A	189	GLN
3	A	191	GLN
3	A	250	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	701	3	14,14,15	0.77	0	17,19,21	1.28	2 (11%)

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	701	3	-	2/6/23/26	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	701	NAG	C1-O5-C5	2.39	115.43	112.19
4	A	701	NAG	C3-C4-C5	2.19	114.15	110.24

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	701	NAG	C3-C2-N2-C7
4	A	701	NAG	C1-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

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	H	222/235 (94%)	0.91	27 (12%) 4 5	24, 50, 100, 126	0
2	L	213/214 (99%)	0.18	12 (5%) 24 29	22, 39, 78, 98	0
3	A	267/274 (97%)	0.53	37 (13%) 2 3	21, 36, 83, 101	0
All	All	702/723 (97%)	0.55	76 (10%) 5 7	21, 41, 93, 126	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	190	GLY	14.1
1	H	189	LEU	13.7
1	H	133	GLY	13.6
1	H	214	LYS	9.1
1	H	191	THR	8.8
1	H	187	SER	8.7
3	A	43	VAL	7.9
3	A	294	PHE	7.6
3	A	45	SER	7.1
1	H	159	LEU	7.0
1	H	207	VAL	6.4
1	H	210	ARG	6.0
3	A	275	ASP	6.0
3	A	46	SER	6.0
1	H	186	SER	5.8
3	A	305	CYS	5.7
3	A	288	ILE	5.7
2	L	122	ASP	5.6
3	A	289	PRO	5.1
3	A	306	PRO	5.0
3	A	48	THR	5.0
2	L	182	SER	5.0
3	A	309	VAL	5.0

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Mol	Chain	Res	Type	RSRZ
3	A	307	LYS	4.9
3	A	308	TYR	4.7
1	H	184	VAL	4.6
1	H	211	VAL	4.6
1	H	198	VAL	4.5
3	A	304	ALA	4.5
3	A	297	VAL	4.4
1	H	212	GLU	4.4
1	H	208	ASP	4.3
1	H	134	GLY	4.1
3	A	293	PRO	4.0
1	H	196	CYS	3.9
1	H	213	PRO	3.7
3	A	291	ASP	3.7
3	A	296	ASN	3.5
1	H	64	LYS	3.4
3	A	279	SER	3.3
3	A	300	ILE	3.3
2	L	184	ALA	3.3
1	H	158	ALA	3.2
3	A	44	GLN	3.2
1	H	185	PRO	3.1
3	A	280	GLU	3.1
1	H	154	TRP	3.1
3	A	287	SER	3.1
2	L	190	LYS	3.0
2	L	212	GLY	3.0
3	A	295	GLN	3.0
2	L	127	SER	2.9
2	L	129	THR	2.9
2	L	185	ASP	2.9
3	A	47	SER	2.8
3	A	290	ASN	2.8
1	H	192	GLN	2.8
1	H	205	THR	2.8
3	A	299	LYS	2.7
3	A	274	ILE	2.7
3	A	49	GLY	2.7
3	A	277	CYS	2.7
3	A	52	CYS	2.6
3	A	278	ILE	2.5
1	H	195	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
3	A	281	CYS	2.4
2	L	189	HIS	2.3
2	L	183	LYS	2.2
1	H	138	LEU	2.2
3	A	276	THR	2.2
3	A	283	THR	2.2
3	A	301	THR	2.2
3	A	51	ILE	2.1
2	L	188	LYS	2.1
2	L	128	GLY	2.1
1	H	160	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](#)

There are no monosaccharides in this entry.

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. 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	B-factors(\AA^2)	Q<0.9
4	NAG	A	701	14/15	0.76	0.29	70,77,82,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.