



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 4, 2017 – 10:55 PM EDT

PDB ID : 4XNM
Title : Antibody Influenza H5 Complex
Authors : Winarski, K.L.; Spiler, B.W.
Deposited on : unknown
Resolution : 2.51 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

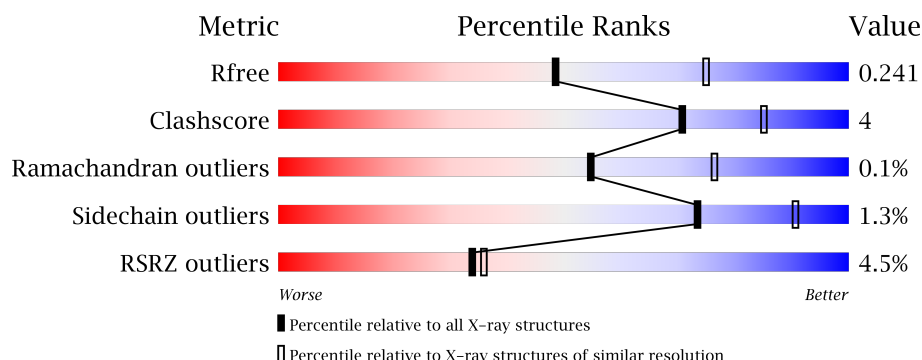
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (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 ≥ 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	208	<div> <div>2%</div> <div>90%</div> <div>10%</div> </div>
1	L	208	<div> <div>84%</div> <div>16%</div> </div>
2	B	227	<div> <div>5%</div> <div>84%</div> <div>11%</div> <div>5%</div> </div>
2	H	227	<div> <div>4%</div> <div>86%</div> <div>9%</div> <div>.</div> </div>
3	C	231	<div> <div>5%</div> <div>75%</div> <div>9%</div> <div>15%</div> </div>

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Mol	Chain	Length	Quality of chain
3	D	231	<div><div></div><div></div><div></div><div></div></div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9769 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 H5.3 FAB Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	208	Total	C	N	O	S	0	0	0
			1569	984	257	322	6			
1	A	208	Total	C	N	O	S	0	0	0
			1569	984	257	322	6			

- Molecule 2 is a protein called H5.3 FAB Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	217	Total	C	N	O	S	0	0	0
			1621	1035	268	314	4			
2	B	216	Total	C	N	O	S	0	0	0
			1617	1033	267	313	4			

- Molecule 3 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	196	Total	C	N	O	S	0	0	0
			1574	1004	272	292	6			
3	D	210	Total	C	N	O	S	0	0	0
			1698	1088	289	315	6			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	38	MET	-	initiating methionine	UNP H8PCX0
C	39	GLY	-	expression tag	UNP H8PCX0
C	40	SER	-	expression tag	UNP H8PCX0
C	41	SER	-	expression tag	UNP H8PCX0
C	42	HIS	-	expression tag	UNP H8PCX0
C	43	HIS	-	expression tag	UNP H8PCX0
C	44	HIS	-	expression tag	UNP H8PCX0
C	45	HIS	-	expression tag	UNP H8PCX0

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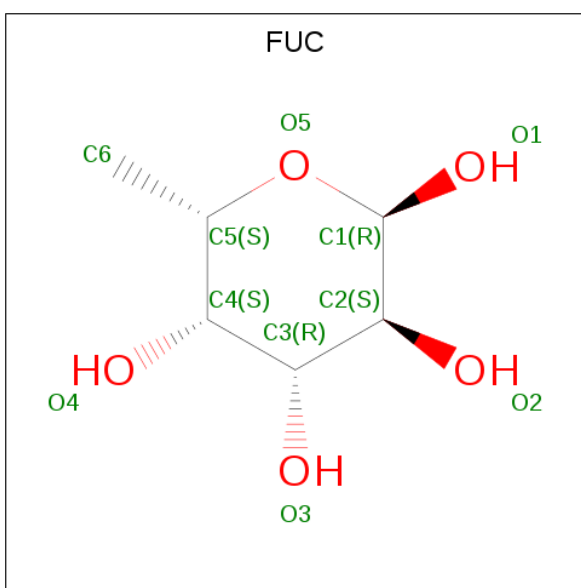
Chain	Residue	Modelled	Actual	Comment	Reference
C	46	HIS	-	expression tag	UNP H8PCX0
C	47	HIS	-	expression tag	UNP H8PCX0
C	48	SER	-	expression tag	UNP H8PCX0
C	49	SER	-	expression tag	UNP H8PCX0
C	50	GLY	-	expression tag	UNP H8PCX0
C	51	LEU	-	expression tag	UNP H8PCX0
C	52	VAL	-	expression tag	UNP H8PCX0
C	53	PRO	-	expression tag	UNP H8PCX0
C	54	ARG	-	expression tag	UNP H8PCX0
C	55	GLY	-	expression tag	UNP H8PCX0
C	56	SER	-	expression tag	UNP H8PCX0
D	38	MET	-	initiating methionine	UNP H8PCX0
D	39	GLY	-	expression tag	UNP H8PCX0
D	40	SER	-	expression tag	UNP H8PCX0
D	41	SER	-	expression tag	UNP H8PCX0
D	42	HIS	-	expression tag	UNP H8PCX0
D	43	HIS	-	expression tag	UNP H8PCX0
D	44	HIS	-	expression tag	UNP H8PCX0
D	45	HIS	-	expression tag	UNP H8PCX0
D	46	HIS	-	expression tag	UNP H8PCX0
D	47	HIS	-	expression tag	UNP H8PCX0
D	48	SER	-	expression tag	UNP H8PCX0
D	49	SER	-	expression tag	UNP H8PCX0
D	50	GLY	-	expression tag	UNP H8PCX0
D	51	LEU	-	expression tag	UNP H8PCX0
D	52	VAL	-	expression tag	UNP H8PCX0
D	53	PRO	-	expression tag	UNP H8PCX0
D	54	ARG	-	expression tag	UNP H8PCX0
D	55	GLY	-	expression tag	UNP H8PCX0
D	56	SER	-	expression tag	UNP H8PCX0

- Molecule 4 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 5 is ALPHA-L-FUCOSE (three-letter code: FUC) (formula: $C_6H_{12}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	L	1	Total	C	O	0	0
			10	6	4		

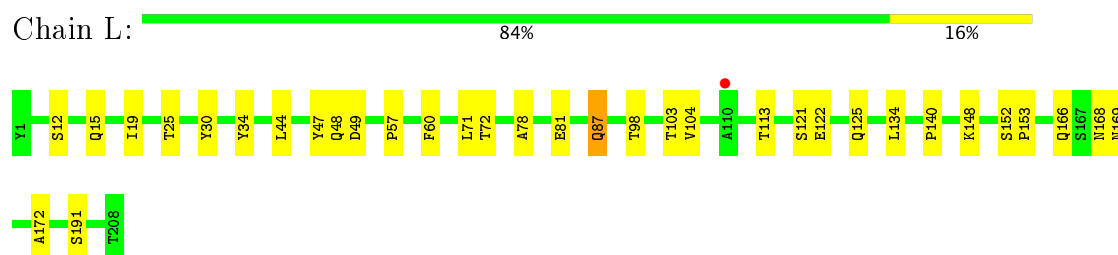
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	L	18	Total 18	O 18	0	0
6	H	9	Total 9	O 9	0	0
6	A	8	Total 8	O 8	0	0
6	B	18	Total 18	O 18	0	0
6	C	20	Total 20	O 20	0	0
6	D	10	Total 10	O 10	0	0

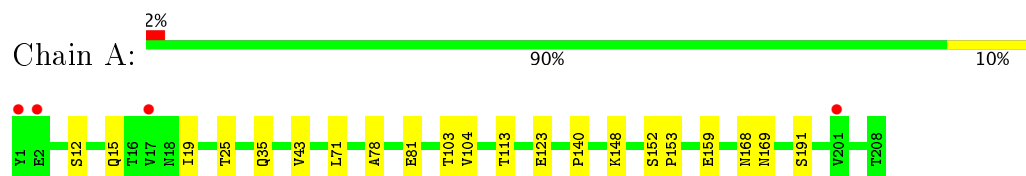
3 Residue-property plots [i](#)

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.

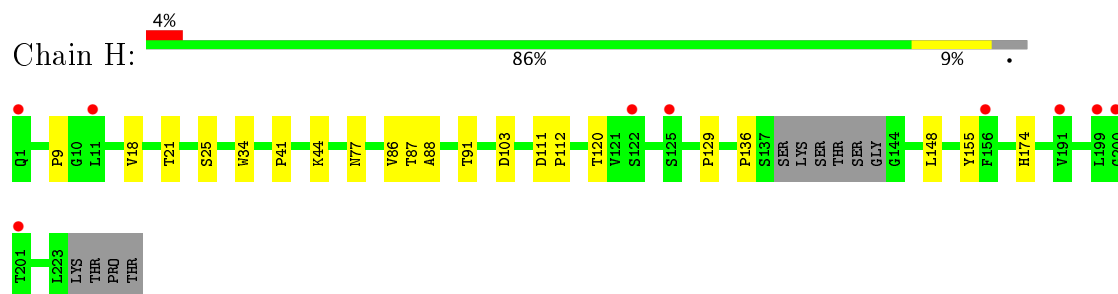
- Molecule 1: H5.3 FAB Light Chain



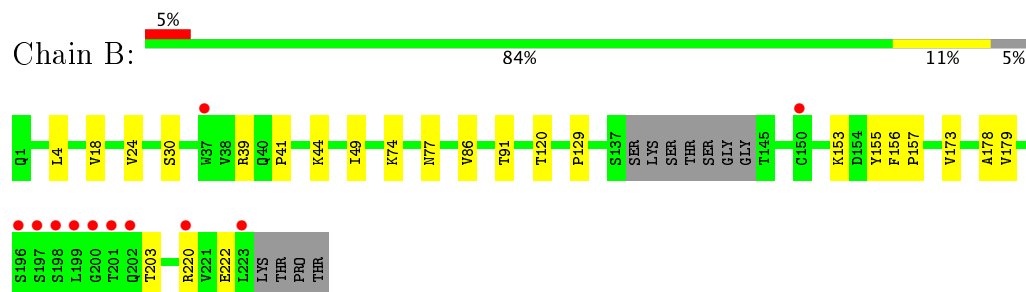
- Molecule 1: H5.3 FAB Light Chain



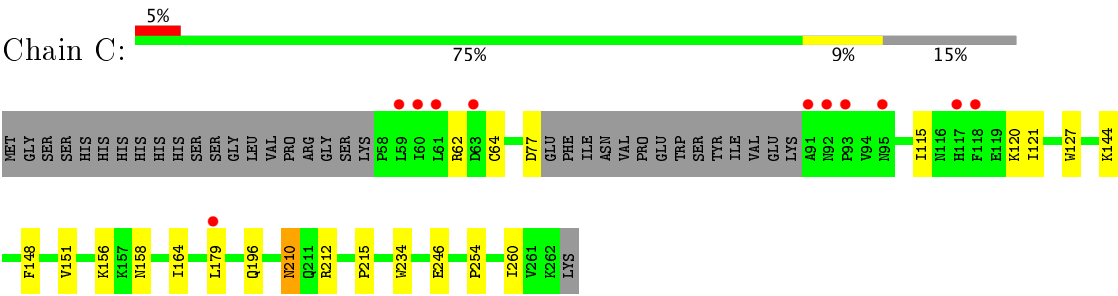
- Molecule 2: H5.3 FAB Heavy Chain



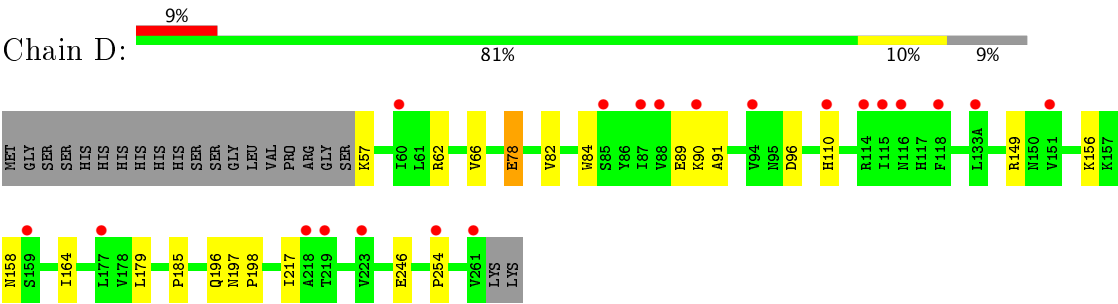
- Molecule 2: H5.3 FAB Heavy Chain



● Molecule 3: Hemagglutinin



● Molecule 3: Hemagglutinin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	250.94Å 51.43Å 127.18Å 90.00° 99.83° 90.00°	Depositor
Resolution (Å)	29.46 – 2.51 29.46 – 2.51	Depositor EDS
% Data completeness (in resolution range)	96.8 (29.46-2.51) 90.0 (29.46-2.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1760)	Depositor
R, R_{free}	0.214 , 0.241 0.212 , 0.241	Depositor DCC
R_{free} test set	1851 reflections (3.70%)	DCC
Wilson B-factor (Å ²)	60.2	Xtriage
Anisotropy	0.475	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9769	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, FUC

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	0.21	0/1609	0.40	0/2201
1	L	0.22	0/1609	0.40	0/2201
2	B	0.21	0/1660	0.41	0/2272
2	H	0.21	0/1664	0.41	0/2277
3	C	0.21	0/1616	0.38	0/2196
3	D	0.21	0/1746	0.39	0/2377
All	All	0.21	0/9904	0.40	0/13524

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	1569	0	1513	13	0
1	L	1569	0	1512	21	0
2	B	1617	0	1600	15	0
2	H	1621	0	1603	9	0
3	C	1574	0	1536	15	0
3	D	1698	0	1653	12	0
4	L	28	0	24	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	L	10	0	10	0	0
6	A	8	0	0	1	0
6	B	18	0	0	1	0
6	C	20	0	0	4	0
6	D	10	0	0	2	0
6	H	9	0	0	0	0
6	L	18	0	0	5	0
All	All	9769	0	9451	78	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 4.

All (78) 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:L:34:TYR:HE2	1:L:87:GLN:HG2	1.36	0.87
1:L:98:THR:OG1	6:L:401:HOH:O	1.97	0.83
1:A:159:GLU:HB3	2:B:179:VAL:HG21	1.66	0.78
1:L:49:ASP:O	6:L:402:HOH:O	2.01	0.78
3:D:96:ASP:OD2	6:D:301:HOH:O	2.02	0.77
3:C:127:TRP:O	6:C:301:HOH:O	2.04	0.76
2:H:129:PRO:HB3	2:H:155:TYR:HB3	1.71	0.72
2:B:129:PRO:HB3	2:B:155:TYR:HB3	1.75	0.69
1:L:78:ALA:HA	1:L:104:VAL:HG21	1.75	0.68
2:B:18:VAL:HG23	2:B:86:VAL:HG11	1.76	0.67
3:D:57:LYS:N	6:D:302:HOH:O	2.28	0.67
1:A:123:GLU:OE2	2:B:153:LYS:NZ	2.28	0.66
1:L:12:SER:HB2	1:L:15:GLN:HG3	1.75	0.66
2:B:220:ARG:NH1	2:B:222:GLU:OE1	2.28	0.66
1:L:122:GLU:OE1	6:L:404:HOH:O	2.14	0.66
3:C:215:PRO:O	6:C:302:HOH:O	2.15	0.64
3:D:185:PRO:HG2	3:D:217:ILE:HG12	1.80	0.63
3:C:179:LEU:HD23	3:C:234:TRP:HB3	1.82	0.62
1:L:81:GLU:HB2	1:L:104:VAL:HG22	1.82	0.62
1:L:103:THR:HG21	1:L:140:PRO:HB3	1.82	0.61
1:A:12:SER:HB2	1:A:15:GLN:HG3	1.83	0.60
2:H:18:VAL:HG23	2:H:86:VAL:HG11	1.84	0.59
1:A:19:ILE:HD12	1:A:71:LEU:HD23	1.85	0.58
2:B:203:THR:HG22	2:B:220:ARG:HH21	1.69	0.58
2:B:30:SER:OG	2:B:74:LYS:NZ	2.28	0.57
1:A:25:THR:HG23	3:D:158:ASN:HA	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:91:THR:HG23	2:H:120:THR:HA	1.88	0.56
1:L:148:LYS:HB2	1:L:191:SER:HB3	1.87	0.56
2:B:178:ALA:HA	2:B:188:LEU:HB3	1.88	0.55
3:C:148:PHE:HB2	3:C:151:VAL:HG12	1.90	0.53
2:B:41:PRO:HG2	2:B:44:LYS:HB2	1.90	0.53
2:B:173:VAL:HG22	2:B:192:VAL:HB	1.92	0.51
1:A:123:GLU:OE1	6:A:302:HOH:O	2.18	0.51
1:L:19:ILE:HD12	1:L:71:LEU:HD23	1.92	0.51
3:D:90:LYS:HD3	3:D:91:ALA:N	2.25	0.51
1:A:148:LYS:HB2	1:A:191:SER:HB3	1.94	0.50
1:A:103:THR:HG21	1:A:140:PRO:HB3	1.93	0.50
1:A:81:GLU:HB2	1:A:104:VAL:HG22	1.93	0.49
3:C:115:ILE:HD12	3:C:260:ILE:HB	1.95	0.48
1:A:78:ALA:HA	1:A:104:VAL:HG21	1.94	0.48
2:H:136:PRO:HD3	2:H:148:LEU:HB3	1.95	0.48
3:D:179:LEU:O	3:D:254:PRO:HB3	2.14	0.48
3:C:156:LYS:HD2	3:C:196:GLN:HB2	1.96	0.48
3:C:179:LEU:O	3:C:254:PRO:HB3	2.14	0.48
2:B:4:LEU:HD22	2:B:24:VAL:HG22	1.96	0.47
3:C:210:ASN:OD1	3:D:149:ARG:NH2	2.47	0.47
3:C:164:ILE:O	3:C:246:GLU:HA	2.14	0.47
1:L:48:GLN:NE2	2:H:103:ASP:OD1	2.46	0.47
1:L:34:TYR:CE2	1:L:87:GLN:HG2	2.29	0.47
3:C:212:ARG:NE	6:C:307:HOH:O	2.38	0.46
1:A:168:ASN:O	1:A:169:ASN:HB2	2.16	0.46
1:L:57:PRO:HG2	1:L:60:PHE:CD1	2.52	0.45
2:H:41:PRO:HG2	2:H:44:LYS:HB2	1.99	0.45
1:L:72:THR:OG1	6:L:406:HOH:O	2.20	0.45
1:L:25:THR:HG23	3:C:158:ASN:HA	1.98	0.45
3:D:62:ARG:N	3:D:78:GLU:OE2	2.41	0.45
1:L:166:GLN:OE1	1:L:172:ALA:HB2	2.17	0.45
2:B:91:THR:HG23	2:B:120:THR:HA	1.98	0.44
2:H:34:TRP:CE2	3:C:144:LYS:HE3	2.53	0.44
1:L:168:ASN:O	1:L:169:ASN:HB2	2.17	0.44
2:H:87:THR:OG1	2:H:88:ALA:N	2.51	0.44
1:L:44:LEU:HD11	1:L:47:TYR:HB3	1.99	0.43
1:A:152:SER:HA	1:A:153:PRO:HD3	1.88	0.43
3:C:120:LYS:O	3:C:121:ILE:HD12	2.19	0.43
1:L:30:TYR:O	6:L:407:HOH:O	2.21	0.43
3:C:62:ARG:O	3:C:64:CYS:N	2.46	0.43
1:L:121:SER:O	1:L:125:GLN:HG3	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:GLN:O	1:A:43:VAL:N	2.51	0.42
2:H:111:ASP:HA	2:H:112:PRO:HA	1.86	0.42
3:C:77:ASP:OD1	6:C:303:HOH:O	2.22	0.42
3:D:156:LYS:HD2	3:D:196:GLN:HB2	2.02	0.42
2:B:156:PHE:HA	2:B:157:PRO:HA	1.82	0.41
3:D:197:ASN:HA	3:D:198:PRO:HD3	1.83	0.41
1:L:152:SER:HA	1:L:153:PRO:HD3	1.93	0.41
2:B:44:LYS:O	6:B:301:HOH:O	2.22	0.41
2:B:39:ARG:HB3	2:B:49:ILE:HD11	2.02	0.40
3:D:164:ILE:O	3:D:246:GLU:HA	2.22	0.40
3:D:66:VAL:HG23	3:D:89:GLU:HB3	2.03	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	206/208 (99%)	203 (98%)	3 (2%)	0	100	100
1	L	206/208 (99%)	204 (99%)	2 (1%)	0	100	100
2	B	212/227 (93%)	207 (98%)	5 (2%)	0	100	100
2	H	213/227 (94%)	207 (97%)	5 (2%)	1 (0%)	32	53
3	C	192/231 (83%)	187 (97%)	5 (3%)	0	100	100
3	D	208/231 (90%)	201 (97%)	7 (3%)	0	100	100
All	All	1237/1332 (93%)	1209 (98%)	27 (2%)	1 (0%)	55	76

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	9	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	179/179 (100%)	178 (99%)	1 (1%)	89	97
1	L	179/179 (100%)	176 (98%)	3 (2%)	66	87
2	B	185/194 (95%)	184 (100%)	1 (0%)	91	97
2	H	185/194 (95%)	181 (98%)	4 (2%)	57	82
3	C	176/208 (85%)	175 (99%)	1 (1%)	89	97
3	D	190/208 (91%)	186 (98%)	4 (2%)	59	83
All	All	1094/1162 (94%)	1080 (99%)	14 (1%)	73	90

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	87	GLN
1	L	113	THR
1	L	134	LEU
2	H	21	THR
2	H	25	SER
2	H	77	ASN
2	H	174	HIS
1	A	113	THR
2	B	77	ASN
3	C	210	ASN
3	D	78	GLU
3	D	82	VAL
3	D	84	TRP
3	D	110	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	64	ASN
1	A	64	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
4	NAG	L	301	1,5,4	14,14,15	0.69	1 (7%)	15,19,21	0.88	1 (6%)
5	FUC	L	302	4	9,10,11	1.37	2 (22%)	13,14,16	1.57	3 (23%)
4	NAG	L	303	4	14,14,15	0.20	0	15,19,21	0.42	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
4	NAG	L	301	1,5,4	-	0/6/23/26	0/1/1/1
5	FUC	L	302	4	-	0/0/17/20	0/1/1/1
4	NAG	L	303	4	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L	301	NAG	O5-C1	-2.32	1.39	1.43
5	L	302	FUC	C2-C3	2.55	1.56	1.52
5	L	302	FUC	C1-C2	3.18	1.59	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	302	FUC	O2-C2-C1	2.29	113.84	109.18
4	L	301	NAG	C1-O5-C5	2.38	115.44	112.17
5	L	302	FUC	C1-O5-C5	2.83	118.64	112.39
5	L	302	FUC	C1-C2-C3	3.15	113.64	109.65

There are no chirality outliers.

There are no torsion outliers.

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 ⓘ

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	208/208 (100%)	0.12	4 (1%) 67 69	64, 92, 117, 136	0
1	L	208/208 (100%)	0.04	1 (0%) 90 91	53, 79, 110, 129	0
2	B	216/227 (95%)	0.17	12 (5%) 25 26	58, 75, 115, 147	0
2	H	217/227 (95%)	0.21	9 (4%) 38 40	55, 85, 119, 142	0
3	C	196/231 (84%)	0.30	11 (5%) 25 26	49, 72, 129, 145	0
3	D	210/231 (90%)	0.43	20 (9%) 9 8	62, 88, 132, 159	0
All	All	1255/1332 (94%)	0.21	57 (4%) 34 36	49, 83, 123, 159	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	114	ARG	7.7
3	D	116	ASN	5.6
2	B	201	THR	4.6
3	C	59	LEU	4.6
3	C	91	ALA	3.8
3	C	61	LEU	3.7
3	C	93	PRO	3.6
2	B	196	SER	3.6
3	D	110	HIS	3.3
2	H	201	THR	3.2
2	H	200	GLY	3.2
2	H	1	GLN	3.1
2	H	199	LEU	3.1
2	H	11	LEU	3.1
1	A	201	VAL	3.1
1	A	17	VAL	2.9
3	D	218	ALA	2.9
3	D	261	VAL	2.9
3	C	118	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
3	D	159	SER	2.8
3	C	63	ASP	2.8
3	C	60	ILE	2.7
2	H	125	SER	2.7
2	H	122	SER	2.6
3	C	179	LEU	2.5
3	C	117	HIS	2.5
2	B	37	TRP	2.5
3	D	94	VAL	2.5
2	B	220	ARG	2.5
3	D	85	SER	2.4
3	C	92	ASN	2.4
3	D	219	THR	2.4
3	D	88	VAL	2.4
3	D	223	VAL	2.4
2	B	199	LEU	2.4
2	B	202	GLN	2.3
3	D	115	ILE	2.3
3	D	90	LYS	2.2
1	A	1	TYR	2.2
2	B	150	CYS	2.2
2	B	200	GLY	2.2
3	D	133(A)	LEU	2.2
3	D	177	LEU	2.2
1	L	110	ALA	2.2
3	D	60	ILE	2.2
2	B	197	SER	2.1
3	D	151	VAL	2.1
1	A	2	GLU	2.1
2	B	198	SER	2.1
3	C	95	ASN	2.1
2	H	156	PHE	2.1
3	D	118	PHE	2.1
2	B	195	PRO	2.0
2	B	223	LEU	2.0
2	H	191	VAL	2.0
3	D	87	ILE	2.0
3	D	254	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	NAG	L	301	14/15	0.78	0.24	-	81,88,98,100	0
5	FUC	L	302	10/11	0.88	0.26	-	103,107,158,158	0
4	NAG	L	303	14/15	0.93	0.36	-	102,105,107,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.