



wwPDB X-ray Structure Validation Summary Report i

Mar 1, 2014 – 04:06 AM GMT

PDB ID : 4FP8
Title : Crystal structure of broadly neutralizing antibody C05 bound to H3 influenza hemagglutinin, HA1 subunit
Authors : Ekiert, D.C.; Wilson, I.A.
Deposited on : 2012-06-21
Resolution : 2.95 Å (reported)

This is a wwPDB validation summary 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/ValidationPDFNotes.html>

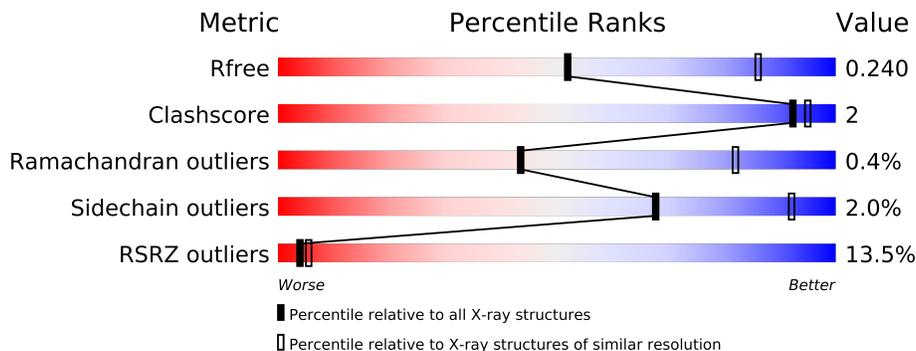
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.95 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}	66092	1424 (2.98-2.90)
Clashscore	79885	1761 (2.98-2.90)
Ramachandran outliers	78287	1708 (2.98-2.90)
Sidechain outliers	78261	1710 (2.98-2.90)
RSRZ outliers	66119	1425 (2.98-2.90)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	278	
1	B	278	
1	C	278	
1	D	278	
2	H	241	
2	I	241	
2	J	241	
2	K	241	
3	L	214	
3	M	214	
3	N	214	
3	O	214	

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 22406 atoms, of which 0 are hydrogen and 0 are deuterium.

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 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	267	2141	1346	375	408	12	0	8	0
1	B	267	2141	1346	375	408	12	0	8	0
1	C	265	2116	1330	369	405	12	0	7	0
1	D	266	2128	1335	374	407	12	0	8	0

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	ALA	-	EXPRESSION TAG	UNP Q91MA7
A	40	ASP	-	EXPRESSION TAG	UNP Q91MA7
A	41	PRO	-	EXPRESSION TAG	UNP Q91MA7
A	42	GLY	-	EXPRESSION TAG	UNP Q91MA7
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
B	39	ALA	-	EXPRESSION TAG	UNP Q91MA7
B	40	ASP	-	EXPRESSION TAG	UNP Q91MA7
B	41	PRO	-	EXPRESSION TAG	UNP Q91MA7
B	42	GLY	-	EXPRESSION TAG	UNP Q91MA7
B	310	GLY	-	EXPRESSION TAG	UNP Q91MA7
B	311	HIS	-	EXPRESSION TAG	UNP Q91MA7
B	312	HIS	-	EXPRESSION TAG	UNP Q91MA7
B	313	HIS	-	EXPRESSION TAG	UNP Q91MA7
B	314	HIS	-	EXPRESSION TAG	UNP Q91MA7
B	315	HIS	-	EXPRESSION TAG	UNP Q91MA7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	316	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	39	ALA	-	EXPRESSION TAG	UNP Q91MA7
C	40	ASP	-	EXPRESSION TAG	UNP Q91MA7
C	41	PRO	-	EXPRESSION TAG	UNP Q91MA7
C	42	GLY	-	EXPRESSION TAG	UNP Q91MA7
C	310	GLY	-	EXPRESSION TAG	UNP Q91MA7
C	311	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	312	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	313	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	314	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	315	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	316	HIS	-	EXPRESSION TAG	UNP Q91MA7
D	39	ALA	-	EXPRESSION TAG	UNP Q91MA7
D	40	ASP	-	EXPRESSION TAG	UNP Q91MA7
D	41	PRO	-	EXPRESSION TAG	UNP Q91MA7
D	42	GLY	-	EXPRESSION TAG	UNP Q91MA7
D	310	GLY	-	EXPRESSION TAG	UNP Q91MA7
D	311	HIS	-	EXPRESSION TAG	UNP Q91MA7
D	312	HIS	-	EXPRESSION TAG	UNP Q91MA7
D	313	HIS	-	EXPRESSION TAG	UNP Q91MA7
D	314	HIS	-	EXPRESSION TAG	UNP Q91MA7
D	315	HIS	-	EXPRESSION TAG	UNP Q91MA7
D	316	HIS	-	EXPRESSION TAG	UNP Q91MA7

- Molecule 2 is a protein called Antibody C05, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	239	Total	C	N	O	S	0	4	0
			1806	1127	306	365	8			
2	I	239	Total	C	N	O	S	0	4	0
			1806	1127	306	365	8			
2	J	239	Total	C	N	O	S	0	4	0
			1806	1127	306	365	8			
2	K	239	Total	C	N	O	S	0	4	0
			1806	1127	306	365	8			

- Molecule 3 is a protein called Antibody C05, light chain.

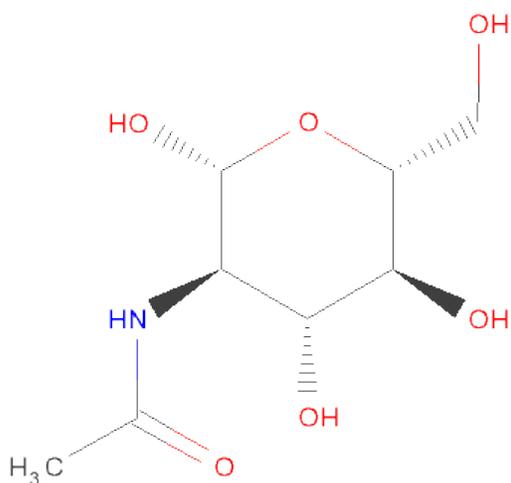
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total	C	N	O	S	0	1	0
			1648	1033	281	330	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	M	213	Total 1648	C 1033	N 281	O 330	S 4	0	1	0
3	N	213	Total 1648	C 1033	N 281	O 330	S 4	0	1	0
3	O	213	Total 1648	C 1033	N 281	O 330	S 4	0	1	0

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



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

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

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

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	39	ALA	-	EXPRESSION TAG	UNP Q91MA7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	40	ASP	-	EXPRESSION TAG	UNP Q91MA7
C	41	PRO	-	EXPRESSION TAG	UNP Q91MA7
C	42	GLY	-	EXPRESSION TAG	UNP Q91MA7
C	310	GLY	-	EXPRESSION TAG	UNP Q91MA7
C	311	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	312	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	313	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	314	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	315	HIS	-	EXPRESSION TAG	UNP Q91MA7
C	316	HIS	-	EXPRESSION TAG	UNP Q91MA7

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

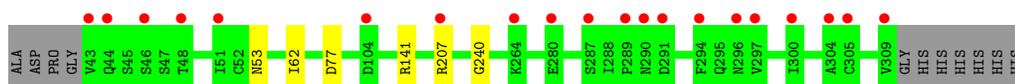
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	J	1	Total Zn 1 1	0	0
6	H	1	Total Zn 1 1	0	0
6	N	2	Total Zn 2 2	0	0
6	O	1	Total Zn 1 1	0	0
6	L	2	Total Zn 2 2	0	0
6	M	1	Total Zn 1 1	0	0

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.

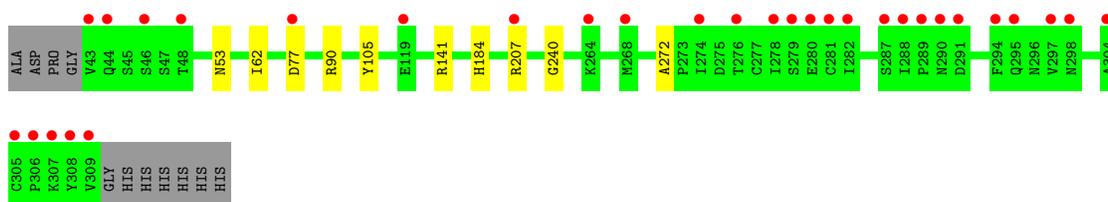
- Molecule 1: Hemagglutinin HA1 chain

Chain A: 



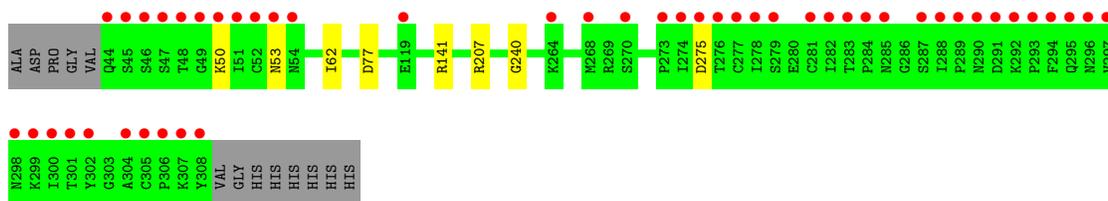
- Molecule 1: Hemagglutinin HA1 chain

Chain B: 



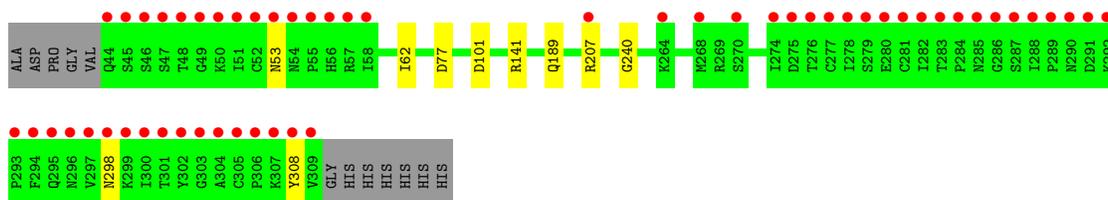
- Molecule 1: Hemagglutinin HA1 chain

Chain C: 



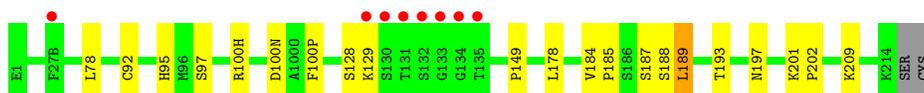
- Molecule 1: Hemagglutinin HA1 chain

Chain D: 



- Molecule 2: Antibody C05, heavy chain

Chain H: 



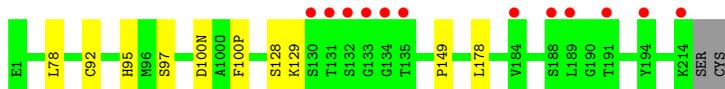
- Molecule 2: Antibody C05, heavy chain

Chain I:



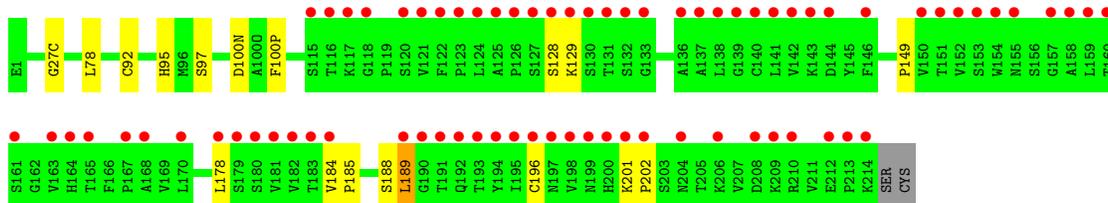
- Molecule 2: Antibody C05, heavy chain

Chain J:



- Molecule 2: Antibody C05, heavy chain

Chain K:



- Molecule 3: Antibody C05, light chain

Chain L:



- Molecule 3: Antibody C05, light chain

Chain M:



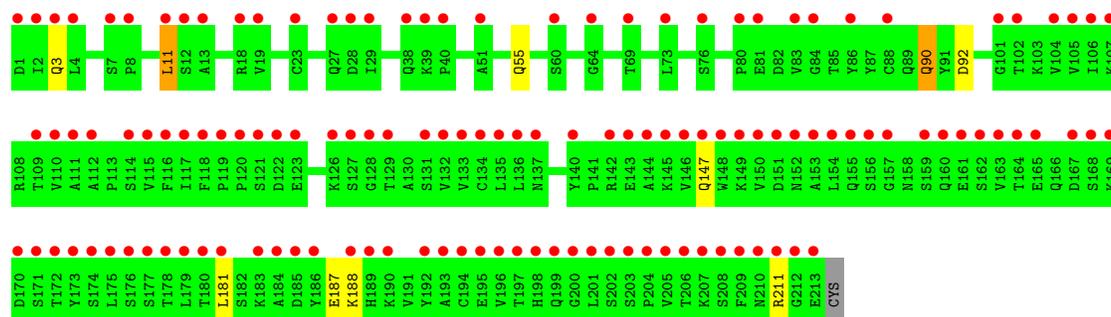
- Molecule 3: Antibody C05, light chain

Chain N:



- Molecule 3: Antibody C05, light chain

Chain O:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.79Å 247.72Å 95.38Å 90.00° 91.33° 90.00°	Depositor
Resolution (Å)	44.49 – 2.95 44.49 – 2.95	Depositor EDS
% Data completeness (in resolution range)	95.1 (44.49-2.95) 95.0 (44.49-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.64 (at 2.96Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, R_{free}	0.212 , 0.241 0.210 , 0.240	Depositor DCC
R_{free} test set	4190 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	56.8	Xtriage
Anisotropy	0.240	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 26.9	EDS
Estimated twinning fraction	0.008 for l,k,-h 0.028 for h,-k,-l 0.021 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 84219 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	22406	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: ZN, PCA, 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	0.30	0/2195	0.58	0/2990
1	B	0.29	0/2195	0.57	0/2990
1	C	0.29	0/2170	0.59	0/2956
1	D	0.31	0/2181	0.58	0/2971
2	H	0.31	0/1837	0.58	1/2500 (0.0%)
2	I	0.28	0/1837	0.54	0/2500
2	J	0.28	0/1837	0.54	0/2500
2	K	0.26	0/1837	0.51	0/2500
3	L	0.34	0/1682	0.59	0/2280
3	M	0.32	0/1682	0.57	0/2280
3	N	0.30	0/1682	0.56	0/2280
3	O	0.26	0/1682	0.52	0/2280
All	All	0.30	0/22817	0.56	1/31027 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	100(H)	ARG	NE-CZ-NH2	-5.03	117.78	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the

chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2141	0	0	2	0
1	B	2141	0	0	5	0
1	C	2116	0	0	3	0
1	D	2128	0	0	5	0
2	H	1806	0	6	6	0
2	I	1806	0	18	3	0
2	J	1806	0	18	3	0
2	K	1806	0	6	7	0
3	L	1648	0	5	2	0
3	M	1648	0	5	2	0
3	N	1648	0	5	1	0
3	O	1648	0	5	4	0
4	A	14	0	13	0	0
4	D	14	0	13	0	0
5	C	28	0	25	0	0
6	H	1	0	0	0	0
6	J	1	0	0	0	0
6	L	2	0	0	0	0
6	M	1	0	0	0	0
6	N	2	0	0	0	0
6	O	1	0	0	0	0
All	All	22406	0	119	40	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

The worst 5 of 40 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:77:ASP:OD2	1:A:141:ARG:NH1	2.16	0.78
1:C:77:ASP:OD2	1:C:141:ARG:NH1	2.22	0.72
1:B:77:ASP:OD2	1:B:141:ARG:NH1	2.32	0.62
3:N:50:ASP:OD1	3:N:91:TYR:OH	2.19	0.60
1:D:77:ASP:OD2	1:D:141:ARG:NH1	2.37	0.57

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	273/278 (98%)	266 (97%)	6 (2%)	1 (0%)	43	82
1	B	273/278 (98%)	266 (97%)	6 (2%)	1 (0%)	43	82
1	C	270/278 (97%)	264 (98%)	5 (2%)	1 (0%)	43	82
1	D	272/278 (98%)	265 (97%)	6 (2%)	1 (0%)	43	82
2	H	241/241 (100%)	234 (97%)	5 (2%)	2 (1%)	27	68
2	I	241/241 (100%)	234 (97%)	6 (2%)	1 (0%)	43	82
2	J	241/241 (100%)	235 (98%)	5 (2%)	1 (0%)	43	82
2	K	241/241 (100%)	233 (97%)	6 (2%)	2 (1%)	27	68
3	L	212/214 (99%)	208 (98%)	4 (2%)	0	100	100
3	M	212/214 (99%)	208 (98%)	4 (2%)	0	100	100
3	N	212/214 (99%)	207 (98%)	5 (2%)	0	100	100
3	O	212/214 (99%)	207 (98%)	5 (2%)	0	100	100
All	All	2900/2932 (99%)	2827 (98%)	63 (2%)	10 (0%)	43	85

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	189	LEU
2	K	189	LEU
1	A	62	ILE
1	B	62	ILE
1	C	62	ILE

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	244/244 (100%)	243 (100%)	1 (0%)	95	99
1	B	244/244 (100%)	243 (100%)	1 (0%)	95	99
1	C	241/244 (99%)	240 (100%)	1 (0%)	95	99
1	D	242/244 (99%)	241 (100%)	1 (0%)	95	99
2	H	202/200 (101%)	195 (96%)	7 (4%)	48	84
2	I	202/200 (101%)	199 (98%)	3 (2%)	76	95
2	J	202/200 (101%)	199 (98%)	3 (2%)	76	95
2	K	202/200 (101%)	198 (98%)	4 (2%)	68	92
3	L	187/187 (100%)	180 (96%)	7 (4%)	45	83
3	M	187/187 (100%)	180 (96%)	7 (4%)	45	83
3	N	187/187 (100%)	180 (96%)	7 (4%)	45	83
3	O	187/187 (100%)	180 (96%)	7 (4%)	45	83
All	All	2527/2524 (100%)	2478 (98%)	49 (2%)	68	93

5 of 49 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	M	11	LEU
3	M	188	LYS
3	O	90	GLN
3	M	90	GLN
2	J	78	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA chains in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length

(or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	PCA	H	1	2	8,8,9	6.56	2 (25%)	8,10,12	6.10	5 (62%)
2	PCA	I	1	2	8,8,9	6.72	2 (25%)	8,10,12	6.64	5 (62%)
2	PCA	J	1	2	8,8,9	6.64	3 (37%)	8,10,12	5.80	5 (62%)
2	PCA	K	1	2	8,8,9	6.67	3 (37%)	8,10,12	5.39	3 (37%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PCA	H	1	2	-	0/0/11/13	0/1/1/1
2	PCA	I	1	2	-	0/0/11/13	0/1/1/1
2	PCA	J	1	2	-	0/0/11/13	0/1/1/1
2	PCA	K	1	2	-	0/0/11/13	0/1/1/1

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	1	PCA	O-C	17.90	1.23	1.11
2	K	1	PCA	O-C	17.70	1.23	1.11
2	J	1	PCA	O-C	17.62	1.23	1.11
2	H	1	PCA	O-C	17.46	1.23	1.11
2	K	1	PCA	CD-N	5.85	1.46	1.34

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	1	PCA	CA-N-CD	-15.28	102.43	114.37
2	J	1	PCA	CA-N-CD	-14.87	102.75	114.37
2	H	1	PCA	CA-N-CD	-14.84	102.77	114.37
2	K	1	PCA	CA-N-CD	-14.43	103.09	114.37
2	I	1	PCA	C-CA-N	-9.63	108.54	110.71

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates (i)

2 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	C	701	1,5	12,14,15	0.67	1 (8%)	15,19,21	0.77	0
5	NAG	C	702	5	12,14,15	0.63	0	15,19,21	1.02	1 (6%)

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	C	701	1,5	-	0/6/23/26	0/1/1/1
5	NAG	C	702	5	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	701	NAG	O5-C5	-2.12	1.41	1.45

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	702	NAG	O5-C5-C6	3.42	110.57	106.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry (i)

Of 10 ligands modelled in this entry, 8 are monoatomic - leaving 2 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
4	NAG	A	701	1	12,14,15	0.74	1 (8%)	15,19,21	0.71	0
4	NAG	D	701	1	12,14,15	0.70	1 (8%)	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
4	NAG	A	701	1	-	0/6/23/26	0/1/1/1
4	NAG	D	701	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	701	NAG	O5-C5	-2.21	1.41	1.45
4	D	701	NAG	O5-C5	-2.14	1.41	1.45

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

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 (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	A	267/278 (96%)	0.38	20 (7%) 14 18	30, 54, 125, 181	0
1	B	267/278 (96%)	0.50	31 (11%) 5 7	29, 60, 146, 202	0
1	C	265/278 (95%)	0.91	48 (18%) 2 3	33, 59, 237, 285	0
1	D	266/278 (95%)	1.34	55 (20%) 1 2	31, 62, 281, 438	0
2	H	239/241 (99%)	0.13	8 (3%) 44 54	31, 50, 97, 265	0
2	I	239/241 (99%)	0.26	8 (3%) 44 54	31, 63, 133, 250	0
2	J	239/241 (99%)	0.32	12 (5%) 28 35	35, 68, 131, 246	0
2	K	239/241 (99%)	2.57	74 (30%) 1 1	36, 100, 312, 488	0
3	L	213/214 (99%)	0.02	2 (0%) 81 88	30, 46, 75, 106	0
3	M	213/214 (99%)	0.03	2 (0%) 81 88	27, 46, 86, 147	0
3	N	213/214 (99%)	0.12	1 (0%) 88 94	49, 69, 113, 140	0
3	O	213/214 (99%)	3.39	129 (60%) 0 0	89, 178, 276, 356	0
All	All	2873/2932 (97%)	0.83	390 (13%) 4 5	27, 63, 228, 488	0

The worst 5 of 390 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	K	182	VAL	23.9
2	K	138	LEU	22.8
2	K	139	GLY	17.9
2	K	140	CYS	17.6
2	K	181	VAL	16.6

6.2 Non-standard residues in protein, DNA, RNA chains (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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	PCA	I	1	8/9	0.18	-	61,81,87,88	0
2	PCA	H	1	8/9	0.21	-	79,96,104,105	0
2	PCA	K	1	8/9	0.22	-	122,139,147,148	0
2	PCA	J	1	8/9	0.17	-	86,98,103,106	0

6.3 Carbohydrates

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	NAG	C	701	14/15	0.33	-	109,120,130,136	0
5	NAG	C	702	14/15	0.32	-	120,125,127,128	0

6.4 Ligands

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	ZN	H	301	1/1	0.16	-	134,134,134,134	0
4	NAG	D	701	14/15	0.31	-	100,115,129,135	0
6	ZN	M	301	1/1	0.08	-	94,94,94,94	0
4	NAG	A	701	14/15	0.28	-	122,134,148,154	0
6	ZN	N	302	1/1	0.07	-	104,104,104,104	0
6	ZN	L	302	1/1	0.13	-	72,72,72,72	0
6	ZN	N	301	1/1	0.13	-	102,102,102,102	0
6	ZN	O	301	1/1	0.11	-	133,133,133,133	0
6	ZN	J	301	1/1	0.05	-	143,143,143,143	0
6	ZN	L	301	1/1	0.09	-	106,106,106,106	0

6.5 Other polymers

There are no such residues in this entry.