



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:54 PM GMT

PDB ID : 4N60
Title : Crystal structure of hemagglutinin from an H7N9 influenza virus in complex with LSTc
Authors : Xu, R.; Wilson, I.A.
Deposited on : 2013-10-11
Resolution : 2.90 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

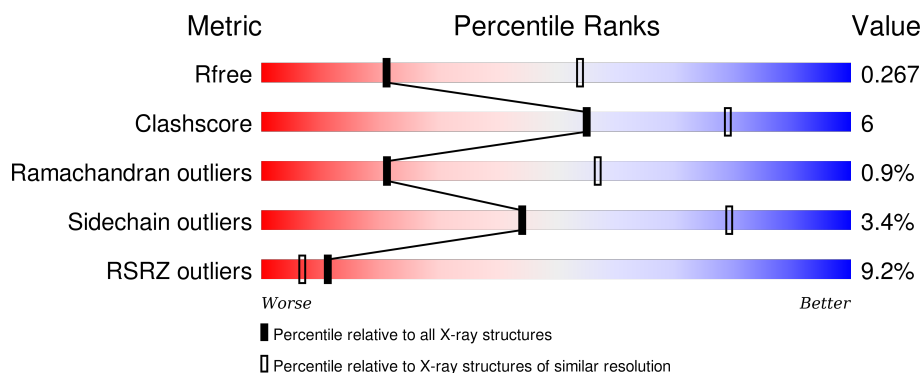
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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. 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	321	<div> <div>2%</div> <div>81%</div> <div>16%</div> <div>..</div> </div>
1	C	321	<div> <div>11%</div> <div>78%</div> <div>19%</div> <div>..</div> </div>
2	B	179	<div> <div>87%</div> <div>6%</div> <div>6%</div> </div>
2	D	179	<div> <div>28%</div> <div>79%</div> <div>15%</div> <div>6%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7770 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	318	Total	C	N	O	S	0	0	0
			2427	1509	438	465	15			
1	C	318	Total	C	N	O	S	0	0	0
			2427	1509	438	465	15			

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	169	Total	C	N	O	S	0	0	0
			1373	845	241	280	7			
2	D	169	Total	C	N	O	S	0	0	0
			1373	845	241	280	7			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	173	SER	-	EXPRESSION TAG	UNP R4NN21
B	174	GLY	-	EXPRESSION TAG	UNP R4NN21
B	175	ARG	-	EXPRESSION TAG	UNP R4NN21
B	176	LEU	-	EXPRESSION TAG	UNP R4NN21
B	177	VAL	-	EXPRESSION TAG	UNP R4NN21
B	178	PRO	-	EXPRESSION TAG	UNP R4NN21
B	179	ARG	-	EXPRESSION TAG	UNP R4NN21
D	173	SER	-	EXPRESSION TAG	UNP R4NN21
D	174	GLY	-	EXPRESSION TAG	UNP R4NN21
D	175	ARG	-	EXPRESSION TAG	UNP R4NN21
D	176	LEU	-	EXPRESSION TAG	UNP R4NN21
D	177	VAL	-	EXPRESSION TAG	UNP R4NN21
D	178	PRO	-	EXPRESSION TAG	UNP R4NN21
D	179	ARG	-	EXPRESSION TAG	UNP R4NN21

- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	2	Total	C	N	O	0	0
			31	17	1	13		
5	C	2	Total	C	N	O	0	0
			31	17	1	13		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	20	Total	O	0	0
			20	20		
6	B	18	Total	O	0	0
			18	18		

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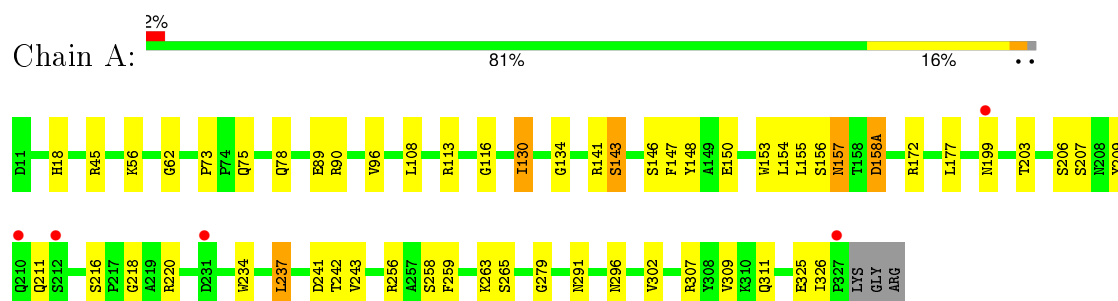
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	3	Total	O	0	0
			3	3		

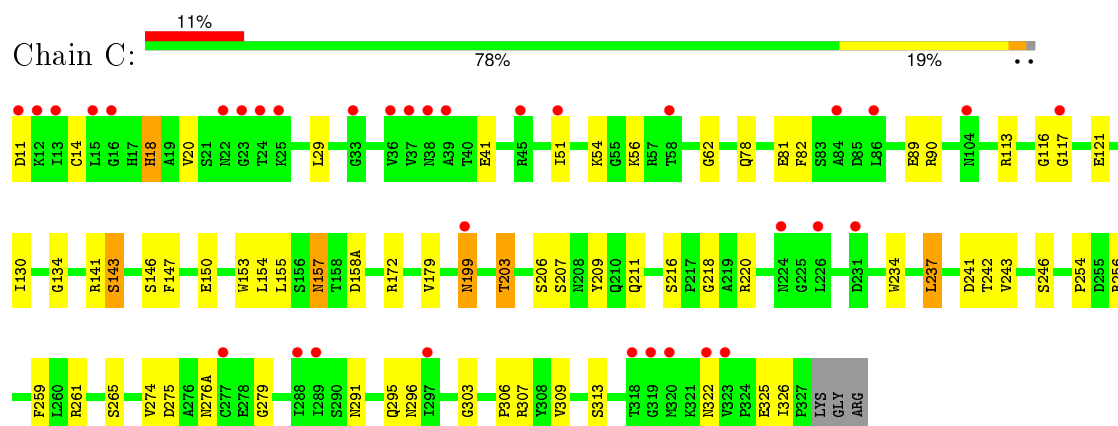
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

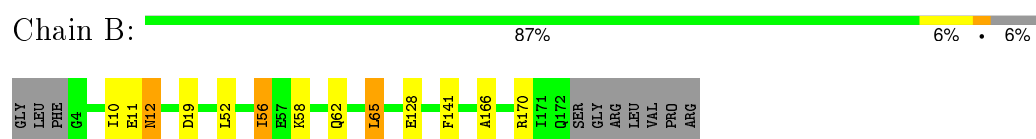
• Molecule 1: Hemagglutinin HA1



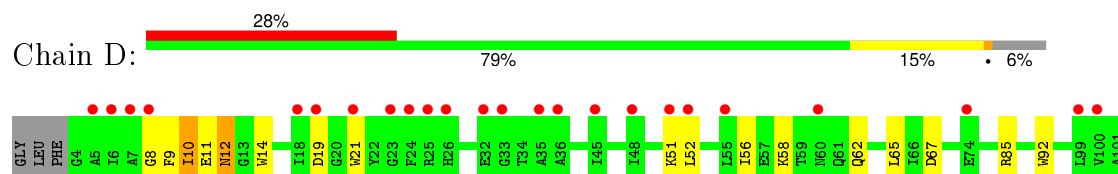
• Molecule 1: Hemagglutinin HA1

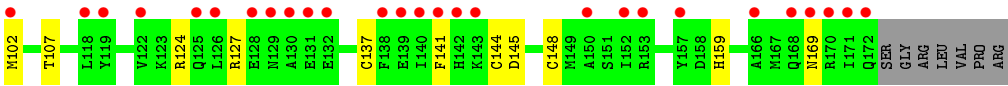


• Molecule 2: Hemagglutinin HA2



• Molecule 2: Hemagglutinin HA2





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	154.77Å 154.77Å 154.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.67 – 2.90 46.67 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (46.67-2.90) 99.6 (46.67-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.223 , 0.262 0.231 , 0.267	Depositor DCC
R_{free} test set	1385 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	71.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 63.4	EDS
Estimated twinning fraction	0.030 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 27536 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7770	wwPDB-VP
Average B, all atoms (Å ²)	123.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.97% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SIA, GAL, BMA, 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	0.25	0/2474	0.44	0/3345
1	C	0.24	0/2474	0.41	0/3345
2	B	0.26	0/1396	0.41	0/1881
2	D	0.22	0/1396	0.39	0/1881
All	All	0.24	0/7740	0.41	0/10452

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	2427	0	2387	34	0
1	C	2427	0	2389	46	0
2	B	1373	0	1273	14	0
2	D	1373	0	1274	21	0
3	A	39	0	34	1	0
4	A	14	0	13	0	0
4	B	14	0	13	0	0
5	A	31	0	26	0	0
5	C	31	0	26	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	20	0	0	1	0
6	B	18	0	0	0	0
6	C	3	0	0	0	0
All	All	7770	0	7435	95	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 6.

All (95) 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:128:GLU:HG3	2:B:170:ARG:HH12	1.29	0.95
1:C:291:ASN:OD1	2:D:58:LYS:NZ	2.13	0.80
2:B:128:GLU:HG3	2:B:170:ARG:NH1	2.00	0.75
2:B:128:GLU:O	2:B:170:ARG:NH1	2.21	0.73
1:C:216:SER:O	1:C:220:ARG:NH2	2.17	0.73
1:A:56:LYS:NZ	1:A:279:GLY:O	2.21	0.73
1:C:296:ASN:ND2	1:C:309:VAL:O	2.24	0.71
1:C:130:ILE:HD11	1:C:154:LEU:HB3	1.73	0.71
1:A:216:SER:O	1:A:220:ARG:NH2	2.20	0.70
1:C:121:GLU:OE2	1:C:172:ARG:NH1	2.25	0.69
1:C:11:ASP:OD2	2:D:144:CYS:N	2.26	0.68
1:A:326:ILE:HG12	2:B:12:ASN:HB3	1.75	0.67
1:C:326:ILE:HG12	2:D:12:ASN:HB3	1.77	0.67
1:A:130:ILE:HD11	1:A:154:LEU:HB3	1.79	0.64
3:A:401:NAG:H61	3:A:402:NAG:HN2	1.62	0.64
1:A:45:ARG:NH1	1:A:296:ASN:OD1	2.31	0.63
1:C:81:GLU:O	1:C:261:ARG:NH2	2.32	0.62
1:C:325:GLU:HA	2:D:12:ASN:HB2	1.81	0.62
1:C:326:ILE:H	2:D:12:ASN:HD22	1.48	0.62
1:C:206:SER:OG	1:C:207:SER:N	2.32	0.60
1:C:303:GLY:HA2	2:D:62:GLN:HG3	1.84	0.59
1:A:150:GLU:OE1	1:A:256:ARG:HD3	2.03	0.59
1:A:237:LEU:HG	1:A:243:VAL:HG13	1.85	0.59
2:D:51:LYS:NZ	2:D:107:THR:OG1	2.30	0.58
1:A:206:SER:OG	1:A:207:SER:N	2.37	0.57
1:C:218:GLY:O	1:C:220:ARG:NH1	2.37	0.57
1:A:291:ASN:HA	2:B:58:LYS:HZ2	1.70	0.57
1:A:172:ARG:HD3	1:A:259:PHE:CZ	2.41	0.56
1:C:295:GLN:HG2	1:C:306:PRO:HG2	1.87	0.56
1:C:199:ASN:N	1:C:199:ASN:OD1	2.32	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:291:ASN:HA	2:D:58:LYS:NZ	2.23	0.54
1:C:237:LEU:HG	1:C:243:VAL:HG13	1.89	0.54
1:C:18:HIS:N	2:D:21:TRP:O	2.36	0.54
1:A:206:SER:HB3	1:A:209:TYR:HB3	1.89	0.54
2:D:67:ASP:OD2	2:D:85:ARG:NH2	2.35	0.54
1:A:291:ASN:HB3	2:B:56:ILE:HG13	1.89	0.54
1:C:146:SER:OG	1:C:147:PHE:N	2.40	0.53
1:C:41:GLU:OE1	1:C:313:SER:OG	2.17	0.53
1:C:203:THR:HG23	1:C:246:SER:HB3	1.90	0.53
1:A:89:GLU:OE2	1:A:113:ARG:NE	2.39	0.52
1:C:89:GLU:OE2	1:C:113:ARG:NH2	2.40	0.52
1:C:291:ASN:HA	2:D:58:LYS:HZ2	1.73	0.52
1:A:134:GLY:HA3	1:A:153:TRP:HB3	1.90	0.52
1:C:155:LEU:HD21	5:C:401:SIA:H111	1.92	0.52
1:A:325:GLU:HA	2:B:12:ASN:HB2	1.93	0.51
1:A:177:LEU:HB3	1:A:258:SER:HB2	1.92	0.51
1:A:291:ASN:OD1	2:B:58:LYS:NZ	2.42	0.50
1:C:29:LEU:HD21	2:D:102:MET:HA	1.94	0.50
1:A:146:SER:OG	1:A:147:PHE:N	2.45	0.49
1:A:291:ASN:HA	2:B:58:LYS:NZ	2.27	0.49
1:C:18:HIS:CG	2:D:21:TRP:HA	2.47	0.49
1:C:275:ASP:OD1	1:C:276(A):ASN:N	2.35	0.49
1:C:56:LYS:NZ	1:C:279:GLY:O	2.47	0.48
1:C:206:SER:HB3	1:C:209:TYR:HB3	1.94	0.48
1:A:307:ARG:NE	6:A:510:HOH:O	2.46	0.48
1:C:130:ILE:HG13	1:C:155:LEU:O	2.14	0.47
1:A:130:ILE:HG13	1:A:155:LEU:O	2.15	0.47
1:C:14:CYS:HA	2:D:137:CYS:HA	1.96	0.47
2:D:127:ARG:HG3	2:D:159:HIS:CG	2.49	0.47
1:C:141:ARG:O	1:C:143:SER:OG	2.29	0.46
1:A:62:GLY:HA2	1:A:90:ARG:HG3	1.97	0.46
1:A:218:GLY:O	1:A:220:ARG:NH1	2.49	0.46
1:C:150:GLU:OE1	1:C:256:ARG:HD3	2.16	0.46
1:C:307:ARG:HG3	2:D:92:TRP:CZ3	2.51	0.45
1:C:116:GLY:HA2	1:C:265:SER:HB3	1.98	0.45
2:D:145:ASP:OD1	2:D:148:CYS:N	2.40	0.45
1:C:172:ARG:HD3	1:C:259:PHE:CZ	2.52	0.44
1:A:130:ILE:HA	1:A:156:SER:O	2.17	0.44
1:A:237:LEU:HD22	1:A:241:ASP:HB3	2.00	0.44
1:C:179:VAL:O	1:C:254:PRO:HB3	2.17	0.44
1:C:134:GLY:HA3	1:C:153:TRP:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:8:GLY:HA2	2:D:14:TRP:HZ2	1.82	0.44
1:A:116:GLY:HA2	1:A:265:SER:HB3	1.99	0.44
1:A:263:LYS:NZ	2:B:62:GLN:OE1	2.43	0.44
1:C:62:GLY:HA2	1:C:90:ARG:HG3	1.99	0.43
1:C:241:ASP:OD1	1:C:242:THR:N	2.43	0.43
1:A:73:PRO:HG3	1:A:148:TYR:CE2	2.53	0.43
2:D:141:PHE:O	2:D:169:ASN:ND2	2.32	0.43
1:A:307:ARG:HH11	1:A:307:ARG:HG2	1.84	0.43
2:B:19:ASP:N	2:B:19:ASP:OD1	2.52	0.43
1:A:141:ARG:O	1:A:143:SER:OG	2.28	0.43
1:C:51:ILE:HB	1:C:274:VAL:HA	2.00	0.42
1:A:302:VAL:HG11	2:B:65:LEU:HD13	2.00	0.42
1:A:241:ASP:OD1	1:A:242:THR:N	2.50	0.41
2:B:65:LEU:HA	2:B:65:LEU:HD12	1.93	0.41
1:A:108:LEU:HD22	1:A:234:TRP:CD1	2.55	0.41
1:A:309:VAL:HB	1:A:311:GLN:OE1	2.20	0.41
2:B:141:PHE:O	2:B:166:ALA:HA	2.20	0.41
1:C:179:VAL:HG22	1:C:234:TRP:HB3	2.03	0.41
1:C:82:PHE:CZ	1:C:117:GLY:HA2	2.55	0.41
2:D:19:ASP:N	2:D:19:ASP:OD1	2.51	0.41
1:C:307:ARG:HH11	1:C:307:ARG:HG2	1.86	0.41
2:D:9:PHE:CD2	2:D:10:ILE:HG13	2.55	0.41
1:C:20:VAL:O	1:C:322:ASN:ND2	2.40	0.40
1:C:54:LYS:HD2	1:C:276(A):ASN:O	2.21	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	316/321 (98%)	294 (93%)	19 (6%)	3 (1%)	21 57

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	316/321 (98%)	294 (93%)	20 (6%)	2 (1%)	30	67
2	B	167/179 (93%)	156 (93%)	9 (5%)	2 (1%)	16	48
2	D	167/179 (93%)	155 (93%)	10 (6%)	2 (1%)	16	48
All	All	966/1000 (97%)	899 (93%)	58 (6%)	9 (1%)	21	57

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	12	ASN
2	D	12	ASN
1	A	157	ASN
1	C	157	ASN
1	A	143	SER
1	C	143	SER
1	A	158(A)	ASP
2	B	10	ILE
2	D	10	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	267/269 (99%)	256 (96%)	11 (4%)	37	73
1	C	267/269 (99%)	259 (97%)	8 (3%)	48	83
2	B	145/153 (95%)	141 (97%)	4 (3%)	51	84
2	D	145/153 (95%)	140 (97%)	5 (3%)	44	79
All	All	824/844 (98%)	796 (97%)	28 (3%)	44	79

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	75	GLN

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Mol	Chain	Res	Type
1	A	78	GLN
1	A	96	VAL
1	A	130	ILE
1	A	157	ASN
1	A	158(A)	ASP
1	A	199	ASN
1	A	203	THR
1	A	211	GLN
1	A	237	LEU
2	B	11	GLU
2	B	52	LEU
2	B	56	ILE
2	B	65	LEU
1	C	18	HIS
1	C	78	GLN
1	C	157	ASN
1	C	158(A)	ASP
1	C	199	ASN
1	C	203	THR
1	C	211	GLN
1	C	237	LEU
2	D	11	GLU
2	D	52	LEU
2	D	56	ILE
2	D	65	LEU
2	D	124	ARG

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

Mol	Chain	Res	Type
1	A	191	GLN
2	D	12	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 ⓘ

7 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	$\# Z > 2$
3	NAG	A	401	1,3	14,14,15	0.65	1 (7%)	15,19,21	0.59	0
3	NAG	A	402	3	14,14,15	0.17	0	15,19,21	0.39	0
3	BMA	A	403	3	11,11,12	0.78	0	14,15,17	0.98	0
5	SIA	A	405	5	16,20,21	0.36	0	18,28,31	0.75	0
5	GAL	A	406	5	11,11,12	0.53	0	14,15,17	0.95	0
5	SIA	C	401	5	16,20,21	0.35	0	18,28,31	0.79	0
5	GAL	C	402	5	11,11,12	0.52	0	14,15,17	0.93	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	401	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	402	3	-	0/6/23/26	0/1/1/1
3	BMA	A	403	3	-	0/2/19/22	0/1/1/1
5	SIA	A	405	5	-	0/14/34/38	1/1/1/1
5	GAL	A	406	5	-	0/2/19/22	0/1/1/1
5	SIA	C	401	5	-	0/14/34/38	1/1/1/1
5	GAL	C	402	5	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	NAG	O5-C1	-2.30	1.39	1.43

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

All (2) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	401	SIA	C2-C3-C4-C5-C6-O6
5	A	405	SIA	C2-C3-C4-C5-C6-O6

3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	NAG	1	0
3	A	402	NAG	1	0
5	C	401	SIA	1	0

5.6 Ligand geometry

2 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	A	404	1	14,14,15	0.22	0	15,19,21	0.43	0
4	NAG	B	201	2	14,14,15	0.39	0	15,19,21	0.43	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	404	1	-	0/6/23/26	0/1/1/1
4	NAG	B	201	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 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 ⓘ

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	318/321 (99%)	-0.10	5 (1%) 74 72	45, 75, 119, 178	0
1	C	318/321 (99%)	0.73	34 (10%) 8 4	87, 135, 232, 299	0
2	B	169/179 (94%)	-0.11	0 100 100	41, 71, 117, 160	0
2	D	169/179 (94%)	1.59	51 (30%) 1 0	99, 213, 283, 337	0
All	All	974/1000 (97%)	0.46	90 (9%) 11 7	41, 106, 250, 337	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	319	GLY	21.0
2	D	7	ALA	10.3
2	D	138	PHE	9.8
2	D	131	GLU	7.8
1	C	13	ILE	7.8
2	D	6	ILE	7.2
2	D	126	LEU	7.2
2	D	130	ALA	7.2
2	D	23	GLY	6.8
2	D	122	VAL	6.5
2	D	36	ALA	6.4
2	D	170	ARG	5.8
1	C	289	ILE	5.6
2	D	129	ASN	5.6
2	D	140	ILE	5.5
1	C	12	LYS	5.5
2	D	172	GLN	5.5
2	D	171	ILE	5.3
1	C	37	VAL	5.2
1	C	58	THR	5.1
1	C	11	ASP	5.1

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Mol	Chain	Res	Type	RSRZ
2	D	24	PHE	5.0
2	D	35	ALA	4.9
1	A	327	PRO	4.7
2	D	166	ALA	4.5
1	C	15	LEU	4.4
2	D	143	LYS	4.4
2	D	157	TYR	4.2
2	D	5	ALA	4.1
1	C	86	LEU	3.8
2	D	119	TYR	3.6
2	D	51	LYS	3.5
2	D	141	PHE	3.4
1	C	323	VAL	3.4
2	D	118	LEU	3.4
1	C	16	GLY	3.3
1	C	33	GLY	3.3
1	C	318	THR	3.3
2	D	48	ILE	3.3
1	C	277	CYS	3.3
2	D	33	GLY	3.3
2	D	142	HIS	3.3
2	D	19	ASP	3.2
1	C	226	LEU	3.2
1	C	22	ASN	3.1
1	C	297	ILE	3.1
2	D	18	ILE	3.1
1	A	199	ASN	3.1
2	D	32	GLU	3.1
2	D	26	HIS	3.1
1	C	288	ILE	3.0
2	D	132	GLU	3.0
1	C	84	ALA	2.9
1	C	38	ASN	2.9
2	D	100	VAL	2.8
1	C	104	ASN	2.8
2	D	25	ARG	2.8
2	D	152	ILE	2.7
1	C	25	LYS	2.7
2	D	74	GLU	2.7
1	A	212	SER	2.7
1	C	51	ILE	2.6
2	D	139	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	45	ARG	2.6
2	D	168	GLN	2.6
2	D	55	LEU	2.5
1	C	23	GLY	2.5
2	D	153	ARG	2.5
1	C	224	ASN	2.4
2	D	45	ILE	2.4
1	A	231	ASP	2.4
2	D	125	GLN	2.4
1	C	199	ASN	2.3
2	D	150	ALA	2.3
1	C	231	ASP	2.3
1	C	320	MET	2.3
2	D	21	TRP	2.3
2	D	102	MET	2.2
2	D	52	LEU	2.2
1	C	24	THR	2.2
1	C	39	ALA	2.2
1	A	210	GLN	2.1
1	C	36	VAL	2.1
1	C	117	GLY	2.1
2	D	8	GLY	2.1
1	C	322	ASN	2.1
2	D	60	ASN	2.1
2	D	128	GLU	2.1
2	D	169	ASN	2.0
2	D	99	LEU	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	SIA	A	405	20/21	0.90	0.24	1.53	89,105,117,117	0
5	SIA	C	401	20/21	0.83	0.36	1.35	131,141,161,161	0
3	NAG	A	402	14/15	0.86	0.28	-	144,151,160,167	0
5	GAL	A	406	11/12	0.85	0.21	-	156,167,171,171	0
5	GAL	C	402	11/12	0.79	0.37	-	199,205,206,206	0
3	NAG	A	401	14/15	0.93	0.18	-	97,111,119,132	0
3	BMA	A	403	11/12	0.66	0.21	-	171,173,174,175	0

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	B	201	14/15	0.89	0.20	-0.11	103,114,119,119	0
4	NAG	A	404	14/15	0.80	0.20	-	146,158,174,177	0

6.5 Other polymers [i](#)

There are no such residues in this entry.